

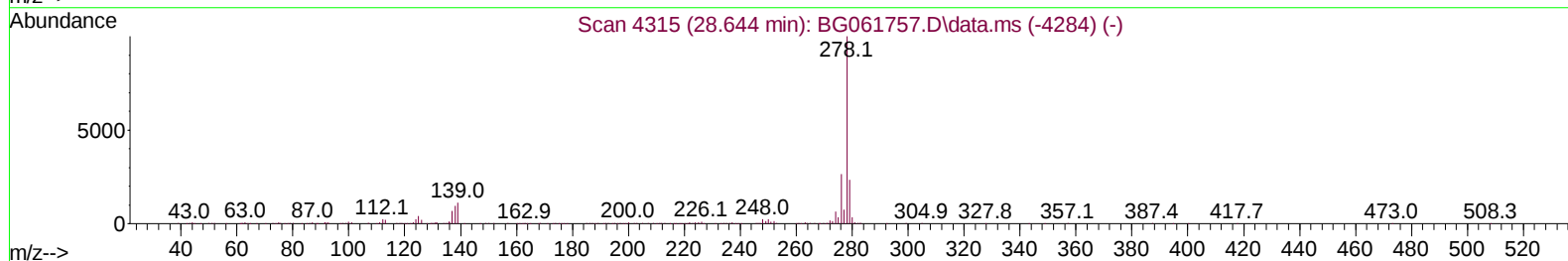
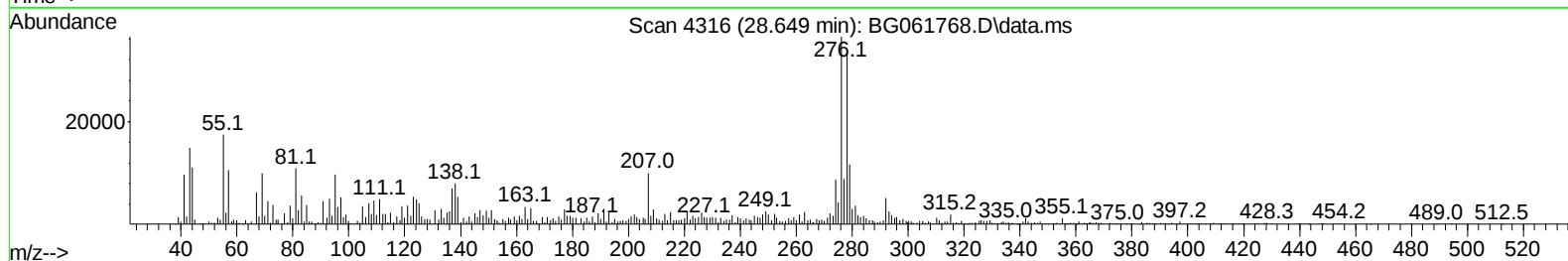
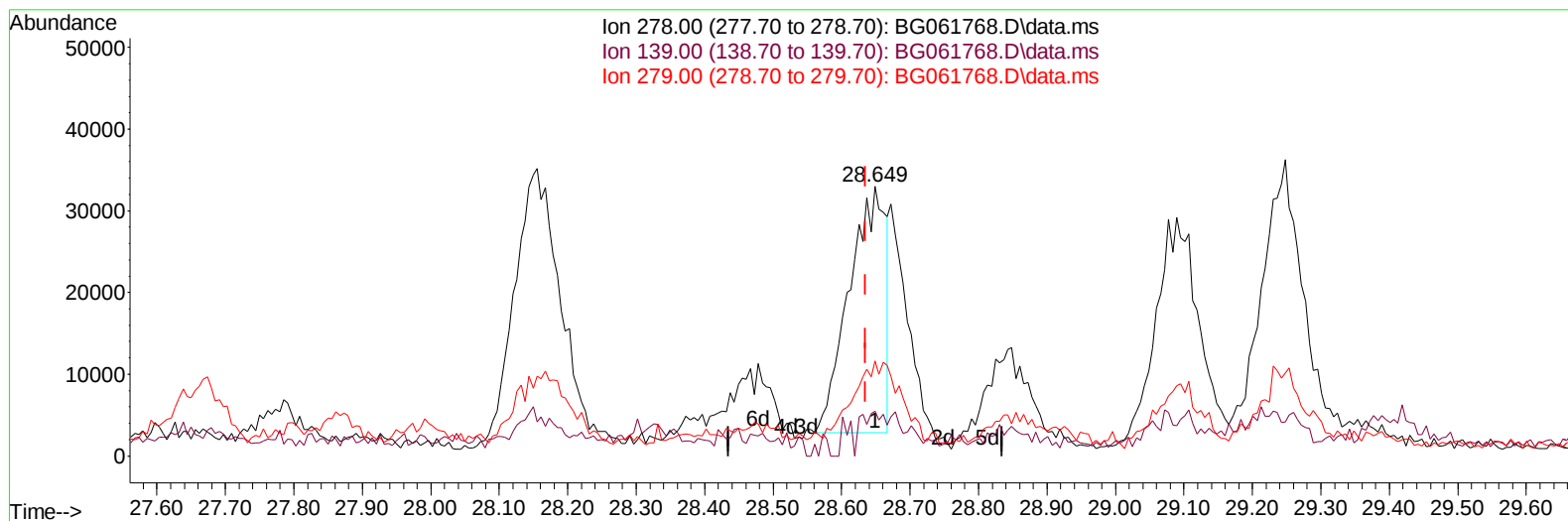
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062724\
 Data File : BG061768.D
 Acq On : 28 Jun 2024 12:09
 Operator : MA/JU
 Sample : P2934-14
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4CQ0

Manual IntegrationsAPPROVED

Quant Time: Jun 28 12:42:45 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 27 12:23:53 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/28/2024
 Supervised By :mohammad ahmed 06/29/2024



TIC: BG061768.D\data.ms

(95) Dibenzo(a,h)anthracene

28.649min (+ 0.014) 4.38 ng/u1

response 110989

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	11.40	16.69#
279.00	23.30	35.37#
0.00	0.00	0.00

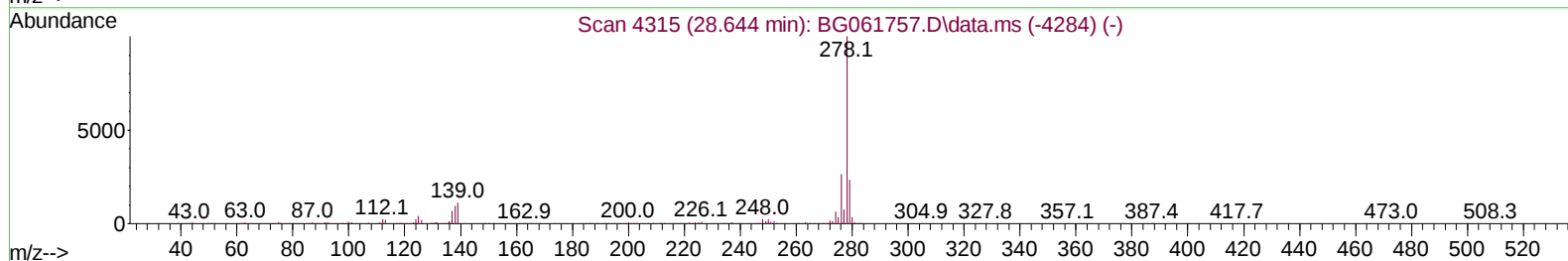
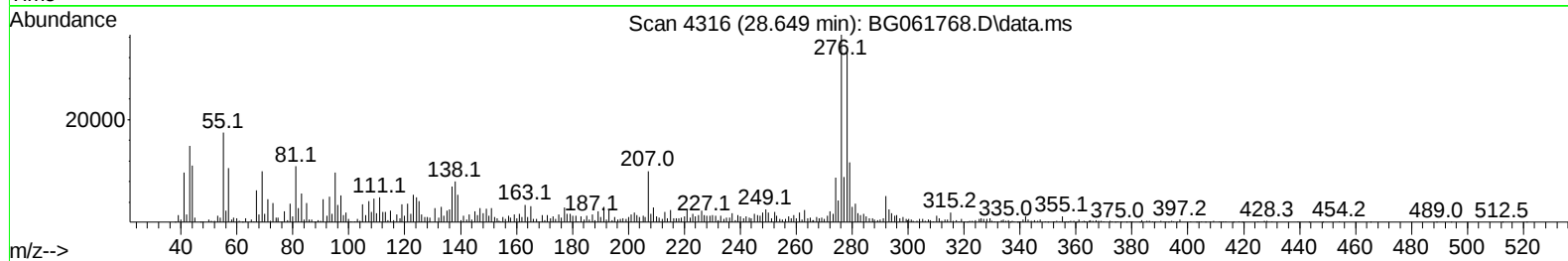
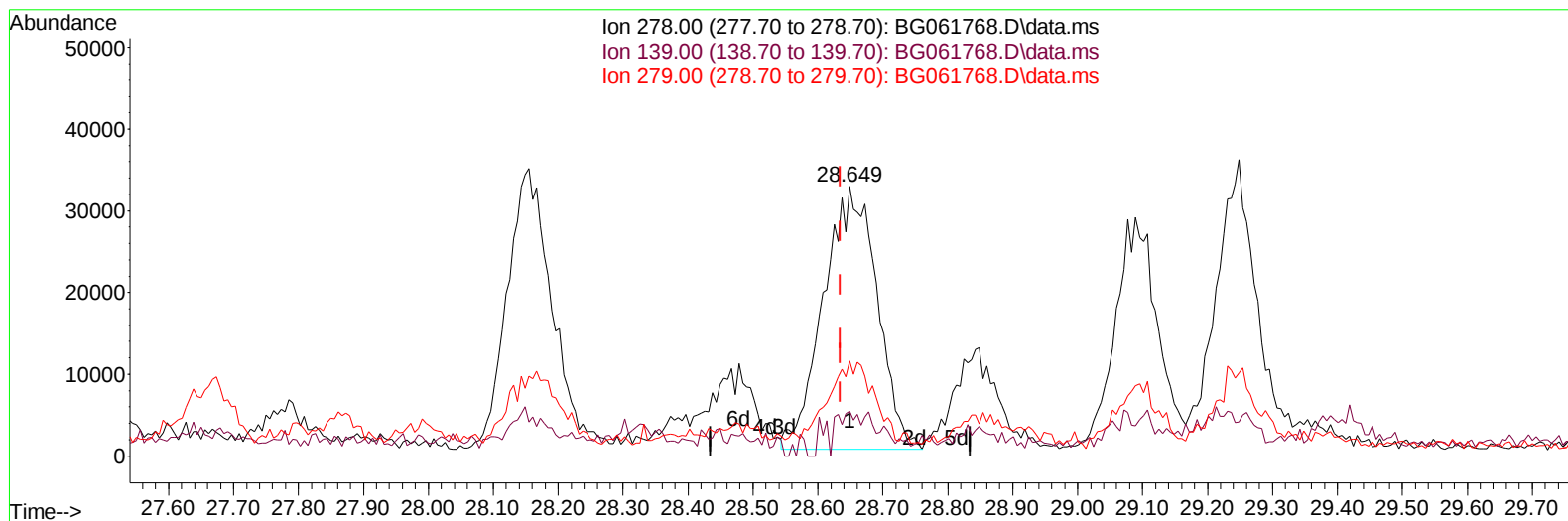
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TIC: BG061768.D\data.ms

(95) Dibenzo(a,h)anthracene

28.649min (+ 0.014) 7.29 ng/ul m

response 184697

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	11.40	16.69#
279.00	23.30	35.37#
0.00	0.00	0.00

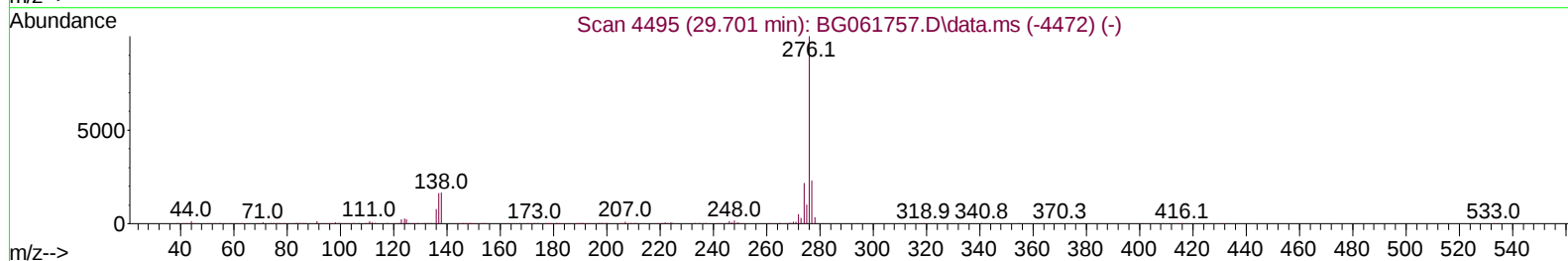
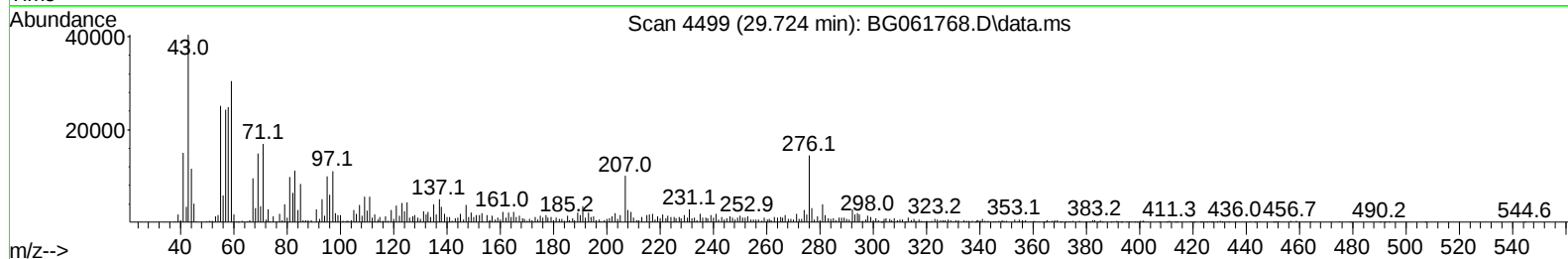
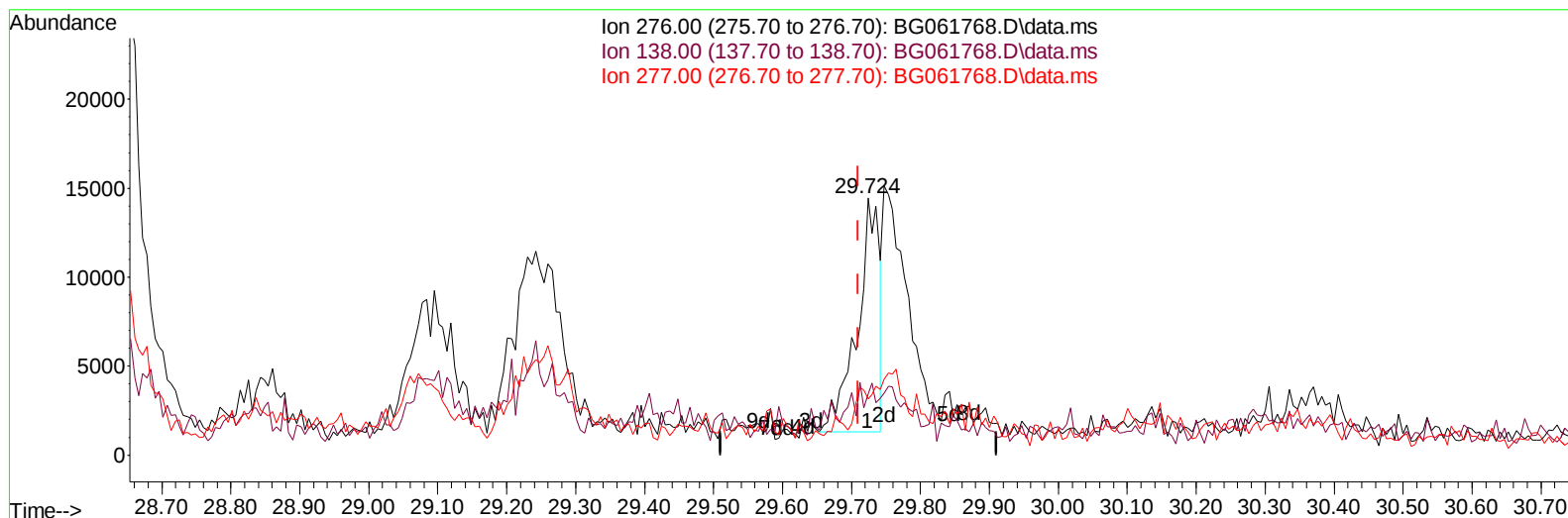
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 Misc :
 ALS Vial : 35 Sample Multiplier: 1

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TIC: BG061768.D\data.ms

(96) Benzo(g,h,i)perylene

29.724min (+ 0.014) 1.19 ng/u1

response 29291

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	15.30	23.77#
277.00	23.20	21.87
0.00	0.00	0.00

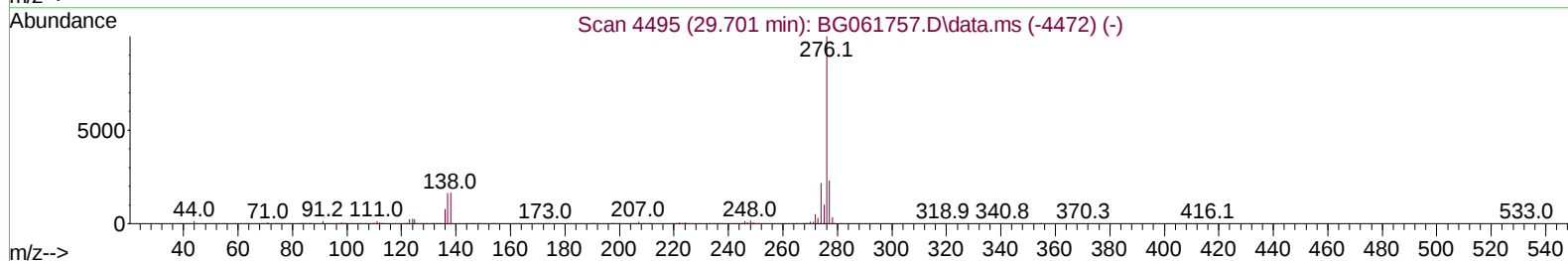
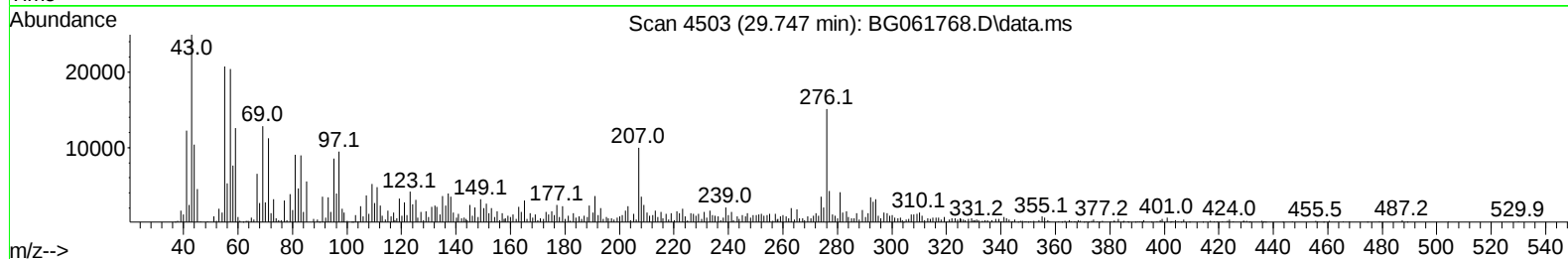
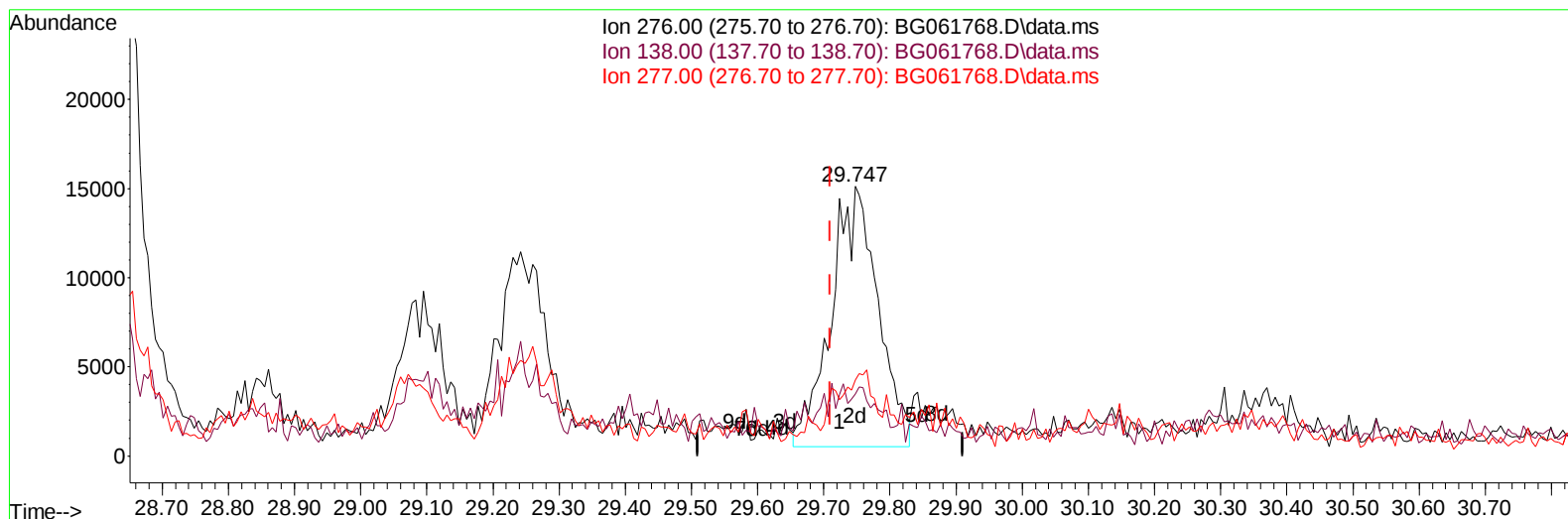
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ALS Vial : 35 Sample Multiplier: 1

Instrument :
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TIC: BG061768.D\data.ms

(96) Benzo(g,h,i)perylene

29.747min (+ 0.037) 2.93 ng/u1 m

response 71956

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	15.30	23.03#
277.00	23.20	28.15#
0.00	0.00	0.00

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 ALS Vial : 35 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Quant Time: Jun 28 12:45:16 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.055	152		85135	20.000	ng/ul	0.00
20) Naphthalene-d8	10.864	136		384057	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.683	164		241674	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.427	188		493140	20.000	ng/ul	0.00
79) Chrysene-d12	21.692	240		365884	20.000	ng/ul	0.00
88) Perylene-d12	24.924	264		422250	20.000	ng/ul	0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.455	96		11225	4.763	ng/uL	0.00
4) Pyridine-d5	3.878	84		28864	4.051	ng/ul	0.00
7) Phenol-d5	7.203	99		258048	27.921	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.380	67		161005	27.945	ng/ul	0.00
11) 2-Chlorophenol-d4	7.579	132		186130	29.348	ng/ul	0.00
15) 4-Methylphenol-d8	8.749	113		192930	27.003	ng/ul	0.00
21) Nitrobenzene-d5	9.219	128		89037	34.144	ng/ul	0.00
24) 2-Nitrophenol-d4	9.941	143		102314	38.739	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.476	165		190428	30.450	ng/ul	0.00
31) 4-Chloroaniline-d4	0.000	131		0d	0.000	ng/ul	0.00
46) Dimethylphthalate-d6	14.084	166		557401	29.971	ng/ul	0.00
49) Acenaphthylene-d8	14.377	160		629417	30.882	ng/ul	0.00
54) 4-Nitrophenol-d4	14.859	143		80448	25.187	ng/ul	0.00
60) Fluorene-d10	15.670	176		469360	28.828	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.781	200		60832	28.848	ng/ul	0.00
73) Anthracene-d10	17.527	188		700079	31.101	ng/ul	0.00
81) Pyrene-d10	19.806	212		600991	29.406	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.700	264		418383	20.746	ng/ul	0.01
Target Compounds							
						Qvalue	
16) Acetophenone	8.878	105		14201	1.207	ng/ul	98
30) Naphthalene	10.917	128		69391	3.229	ng/ul	96
36) 2-Methylnaphthalene	12.521	142		27945	1.843	ng/ul	87
50) Acenaphthylene	14.407	152		58277	2.561	ng/ul#	96
52) Acenaphthene	14.747	153		22619	1.436	ng/ul	93
56) Dibenzofuran	15.076	168		48847	2.215	ng/ul	98
61) Fluorene	15.729	166		29117	1.596	ng/ul#	96
72) Phenanthrene	17.468	178		652458	25.149	ng/ul	99
74) Anthracene	17.562	178		146896	5.537	ng/ul	99
77) Carbazole	17.826	167		48588	2.157	ng/ul#	90
80) Fluoranthene	19.477	202		1192178	50.274	ng/ul	98
82) Pyrene	19.836	202		927390	36.780	ng/ul	98
85) Benzo(a)anthracene	21.675	228		655433	25.520	ng/ul	95
86) Bis(2-ethylhexyl)phtha...	21.586	149		27909	2.068	ng/ul#	96
87) Chrysene	21.739	228		763945	31.435	ng/ul	99
90) Benzo(b)fluoranthene	23.901	252		1071325	42.207	ng/ul	96
91) Benzo(k)fluoranthene	23.966	252		374862	14.360	ng/ul#	97
93) Benzo(a)pyrene	24.771	252		388295	15.890	ng/ul#	92
94) Indeno(1,2,3-cd)pyrene	28.608	276		490823	16.119	ng/ul	99
95) Dibenzo(a,h)anthracene	28.649	278		184697m	7.294	ng/ul	
96) Benzo(g,h,i)perylene	29.747	276		71956m	2.931	ng/ul	

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev	(Min)

(#) = qualifier out of range (m) = manual integration (+) = signals summed								

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