

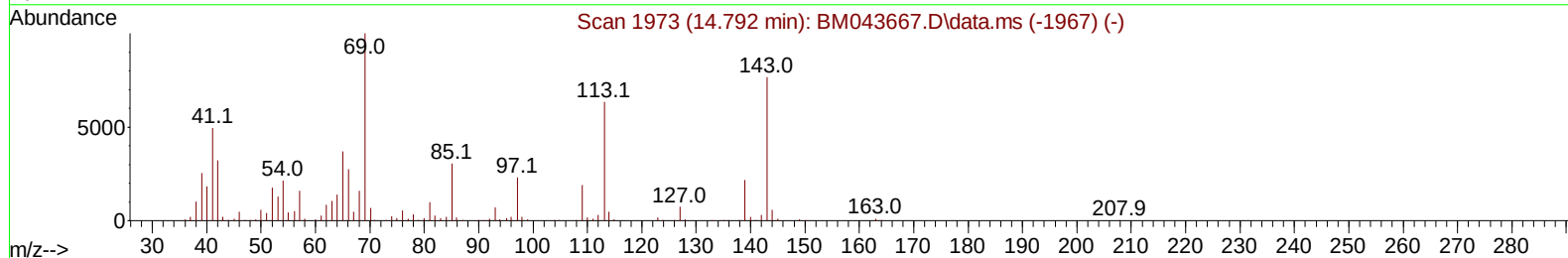
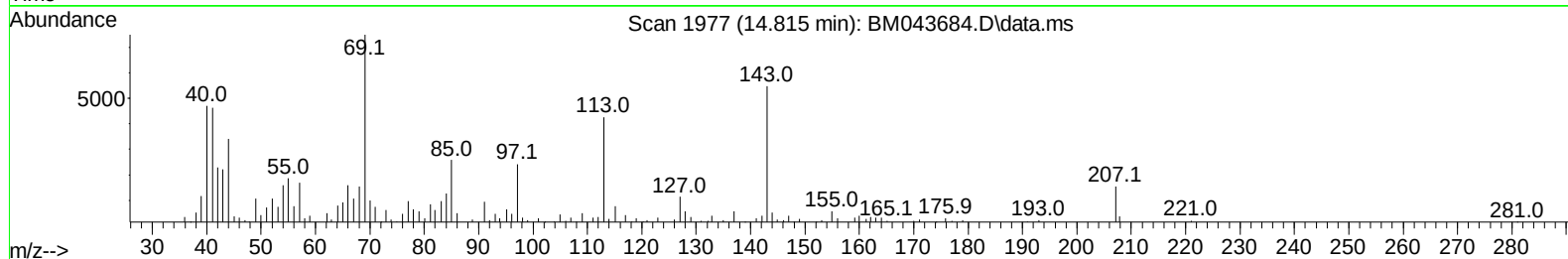
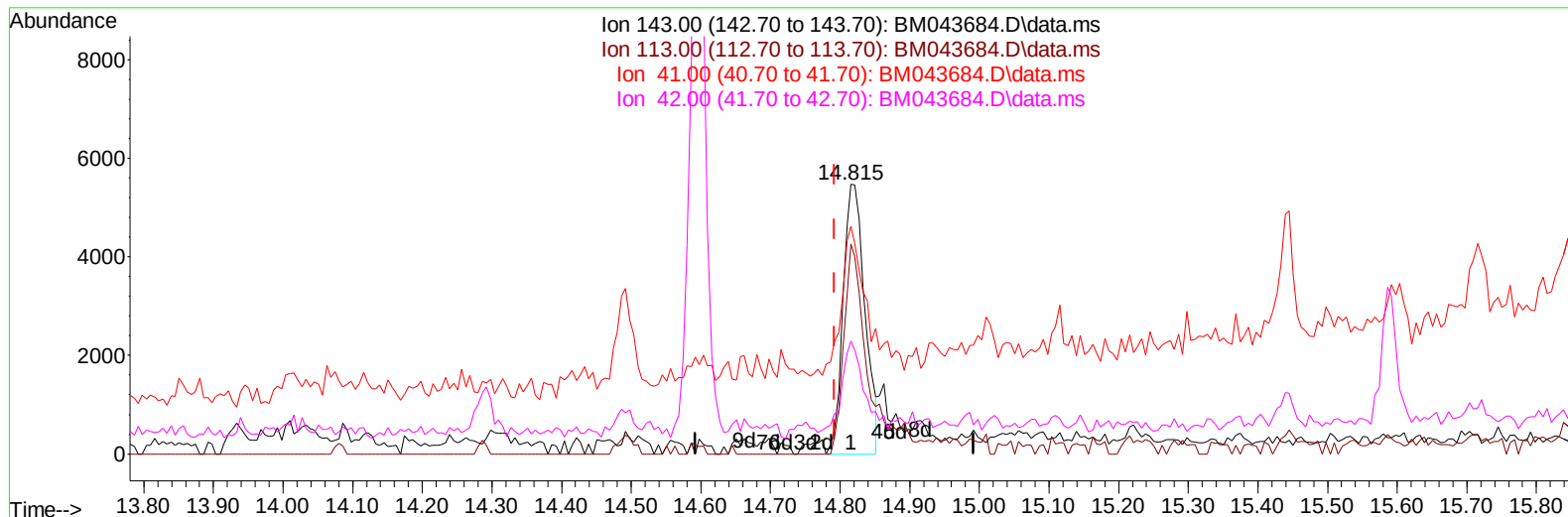
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122923\
 Data File : BM043684.D
 Acq On : 29 Dec 2023 23:18
 Operator : MA/JU
 Sample : 05920-06DL 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GCF57DL

Manual IntegrationsAPPROVED

Quant Time: Dec 30 00:04:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 22:31:57 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 01/01/2024
 Supervised By :mohammad ahmed 01/01/2024



TIC: BM043684.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.815min (+ 0.023) 1.35 ng/ul

response 11938

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.00	77.90
41.00	49.10	84.62#
42.00	33.20	41.83#

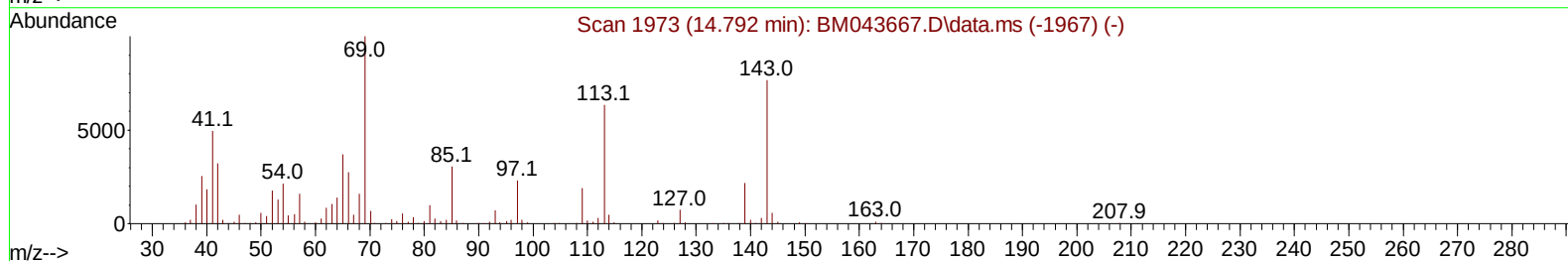
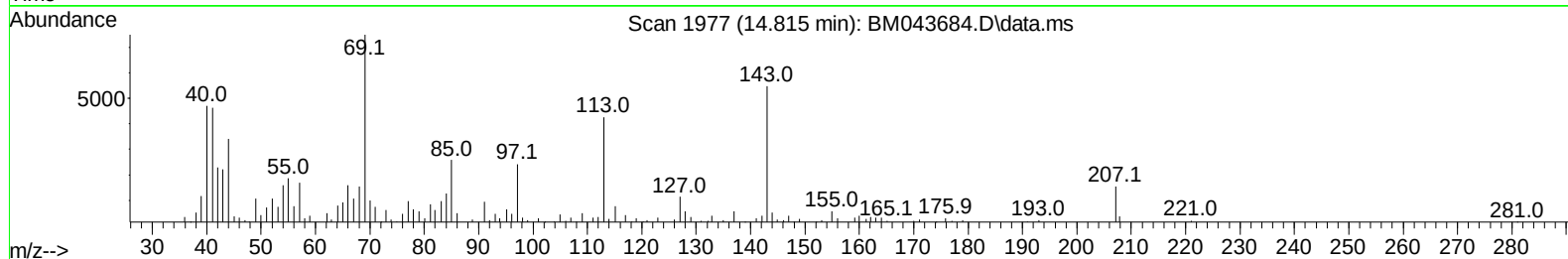
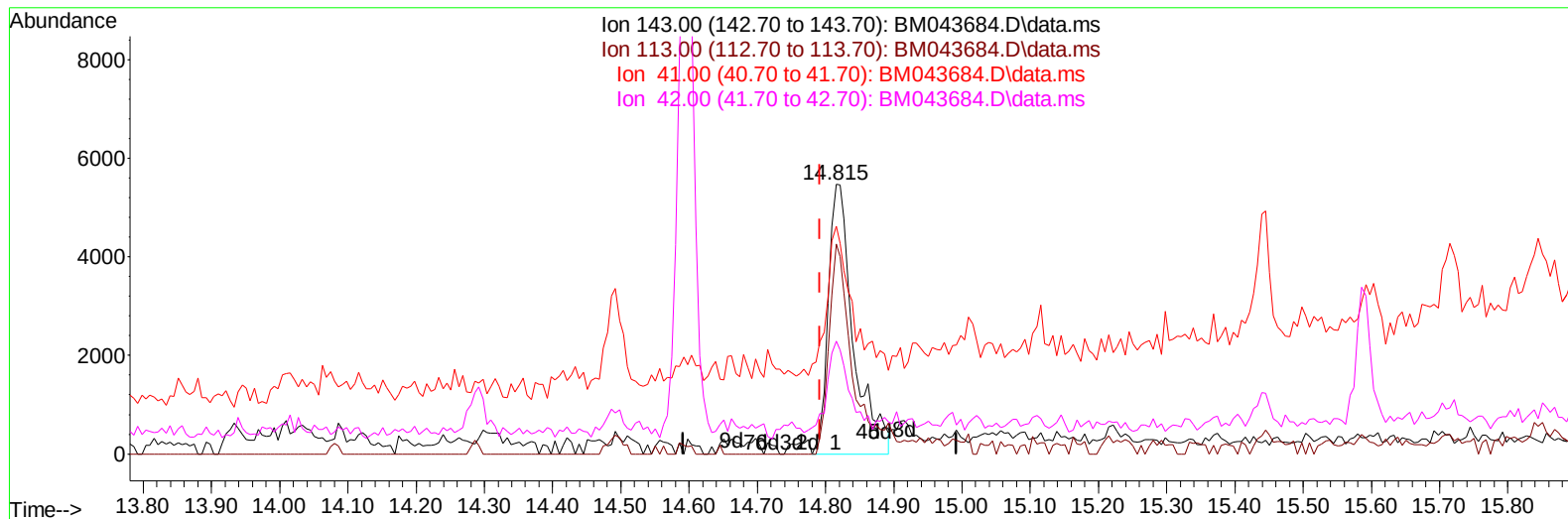
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TIC: BM043684.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.815min (+ 0.023) 1.58 ng/ul m

response 13976

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.00	77.90
41.00	49.10	84.62#
42.00	33.20	41.83#

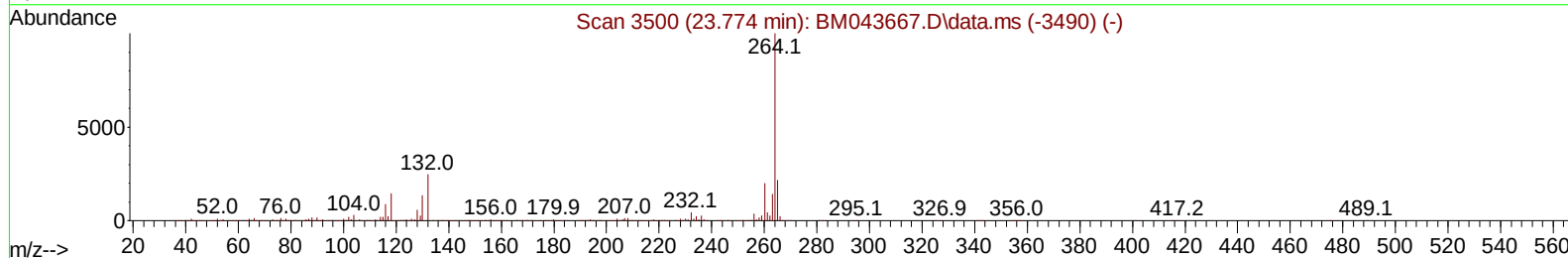
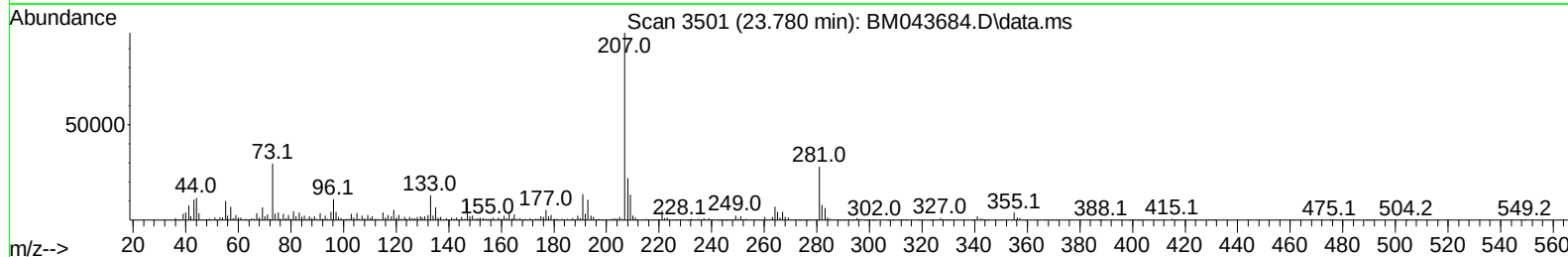
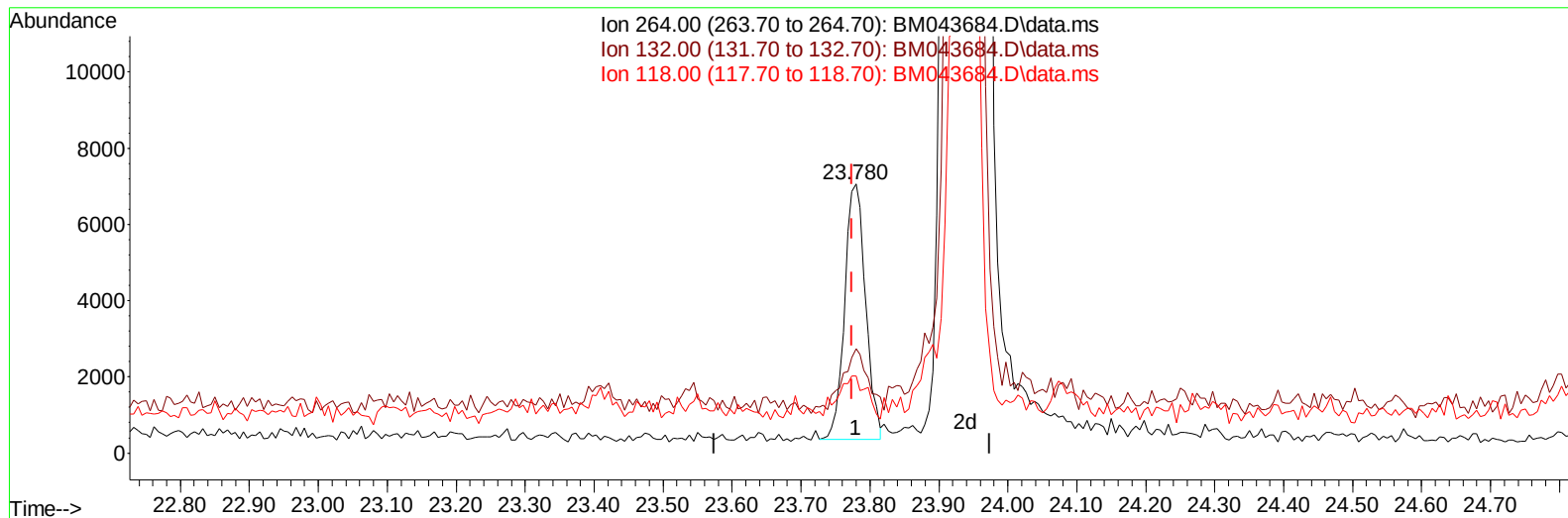
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TIC: BM043684.D\data.ms

(92) Benzo(a)pyrene-d12 (S)

23.780min (+ 0.006) 0.21 ng/u1

response 14081

Ion	Exp%	Act%
264.00	100.00	100.00
132.00	20.20	38.73#
118.00	11.60	28.69#
0.00	0.00	0.00

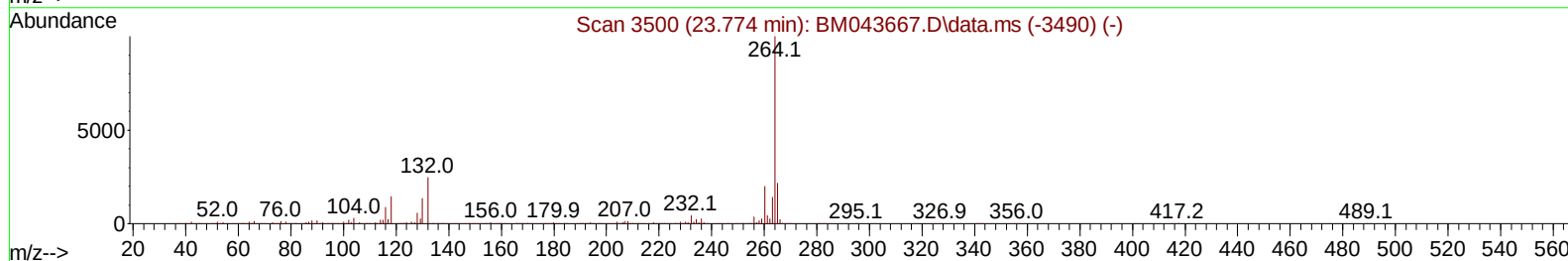
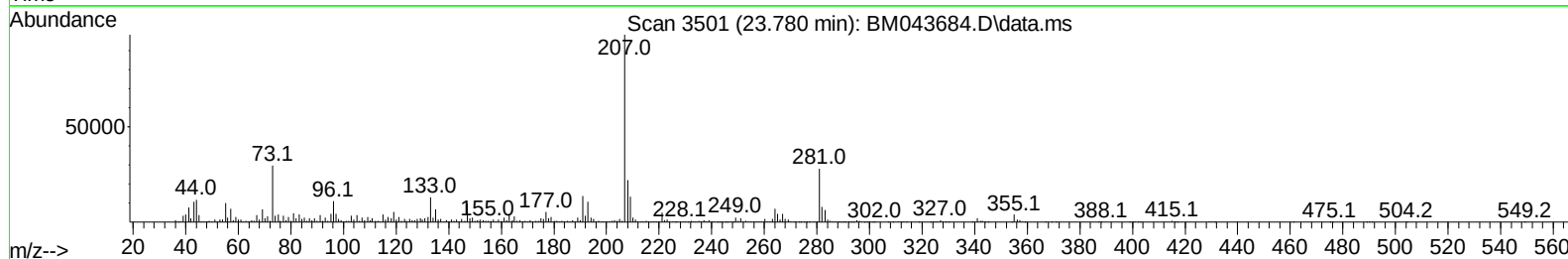
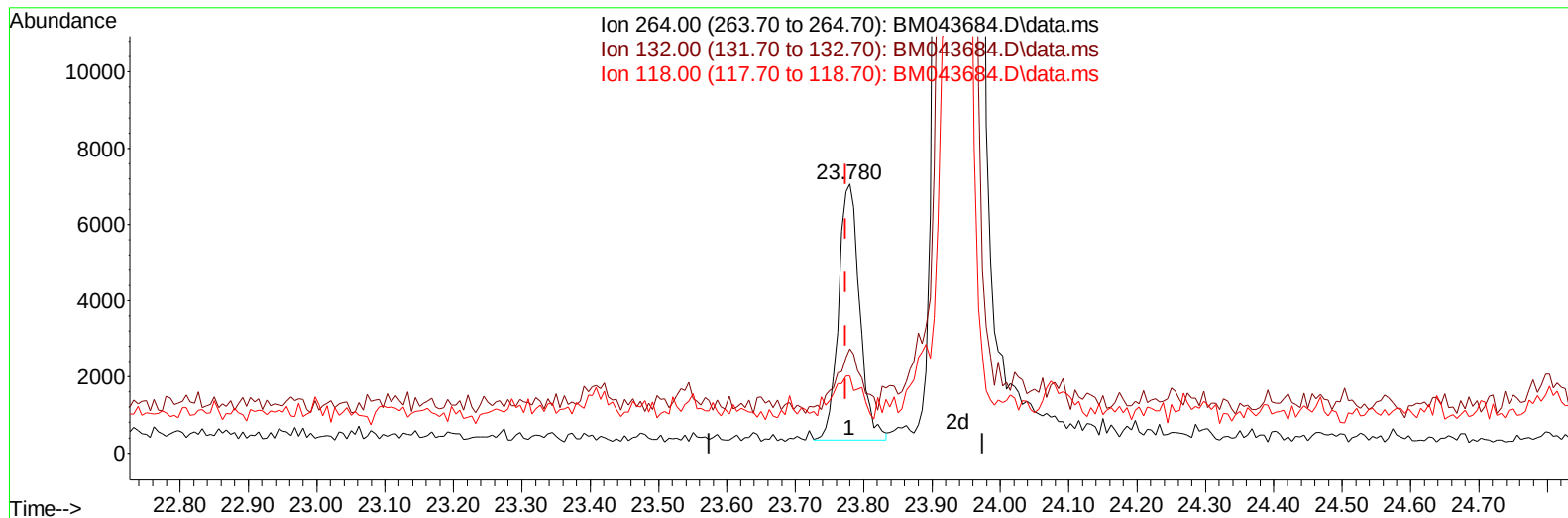
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Instrument :
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TIC: BM043684.D\data.ms

(92) Benzo(a)pyrene-d12 (S)

23.780min (+ 0.006) 0.21 ng/ul m

response 14510

Ion	Exp%	Act%
264.00	100.00	100.00
132.00	20.20	38.73#
118.00	11.60	28.69#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 01/01/2024
 Supervised By :mohammad ahmed 01/01/2024

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	208577	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	907064	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.598	164	518355	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	1115164	20.000	ng/ul	0.00
79) Chrysene-d12	21.521	240	954145	20.000	ng/ul	0.00
88) Perylene-d12	23.938	264	1101449	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	3385	0.527	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.128	99	32685	1.378	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.298	67	42287	2.814	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	43381	2.395	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	34341	1.865	ng/ul	0.00
21) Nitrobenzene-d5	9.133	128	24591	2.802	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143	23704	2.614	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	28363	1.772	ng/ul	0.02
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	14.016	166	130883	2.805	ng/ul	0.00
49) Acenaphthylene-d8	14.292	160	95140	1.719	ng/ul	0.00
54) 4-Nitrophenol-d4	14.815	143	13976m	1.582	ng/ul	0.02
60) Fluorene-d10	15.592	176	122156	3.166	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.439	188	123069	1.968	ng/ul	0.00
81) Pyrene-d10	19.733	212	30914	0.410	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.780	264	14510m	0.213	ng/ul	0.00
Target Compounds						
					Qvalue	
16) Acetophenone	8.798	105	77180	2.738	ng/ul	97
58) 2,3,4,6-Tetrachlorophenol	15.221	232	10605	1.161	ng/ul#	99
71) Pentachlorophenol	16.986	266	403118	44.054	ng/ul	99

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

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