

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121022\
 Data File : BM037937.D
 Acq On : 10 Dec 2022 16:41
 Operator : CG/JU
 Sample : N5896-13ME
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :

BNA_M

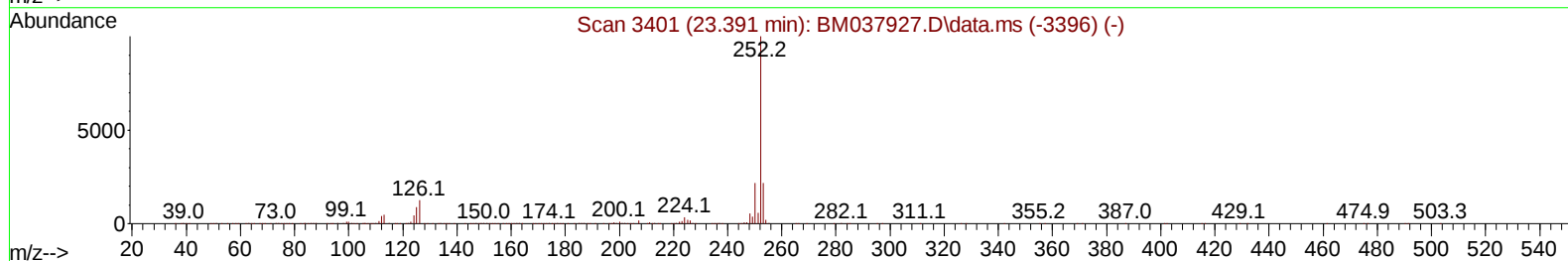
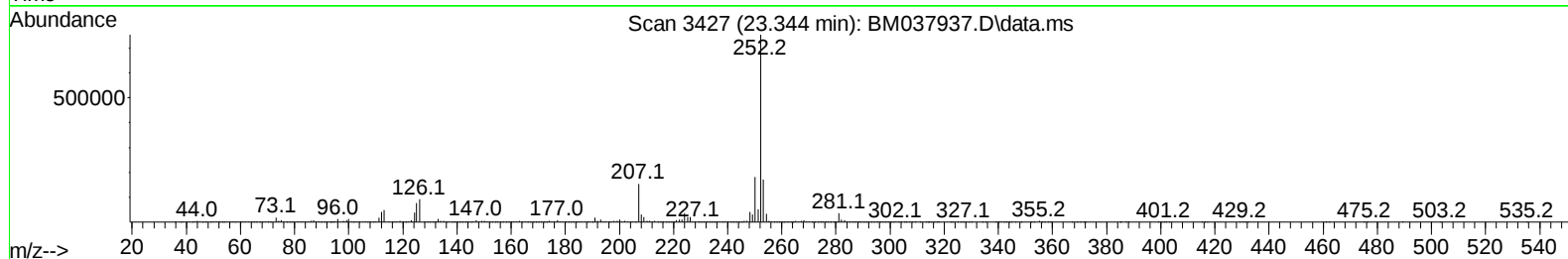
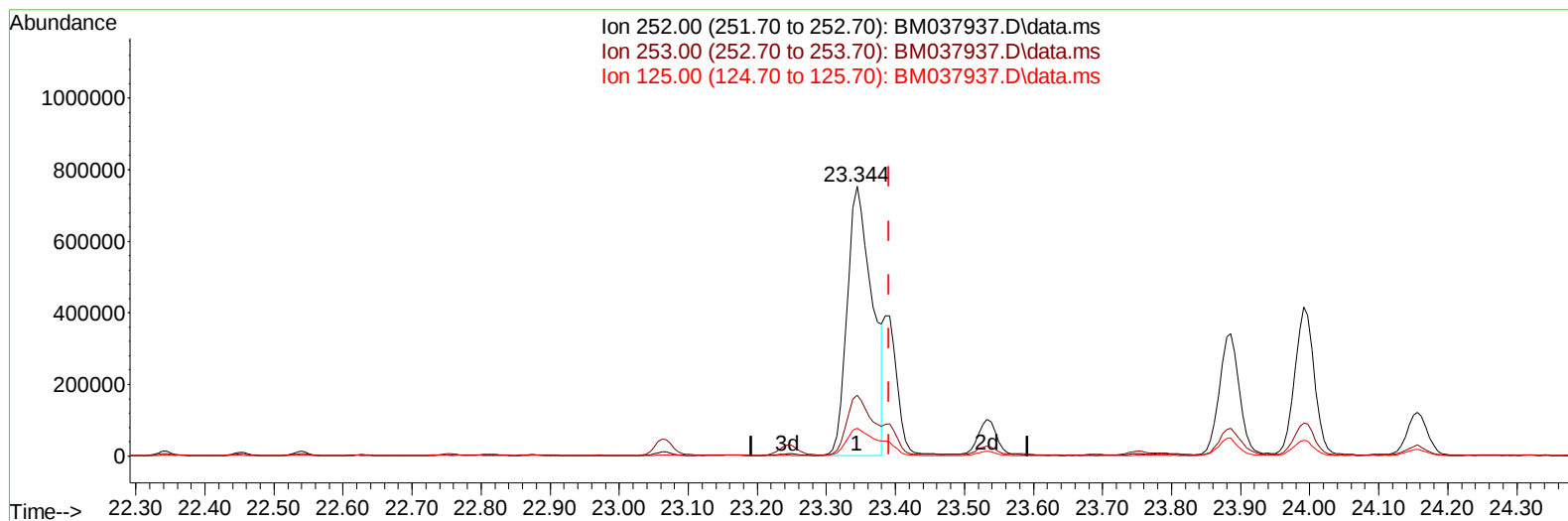
ClientSampleId :

DBXN1ME

Manual IntegrationsAPPROVED

Quant Time: Dec 11 21:27:34 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 09 21:34:57 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/12/2022
 Supervised By :mohammad ahmed 12/12/2022



TIC: BM037937.D\data.ms

(91) Benzo(k)fluoranthene

23.344min (-0.047) 32.01 ng/u1

response 1900536

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.80	22.58
125.00	10.10	10.21
0.00	0.00	0.00

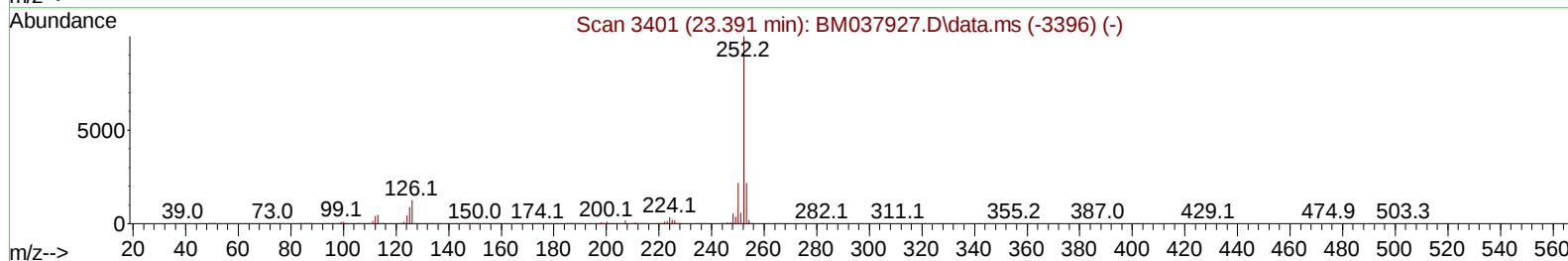
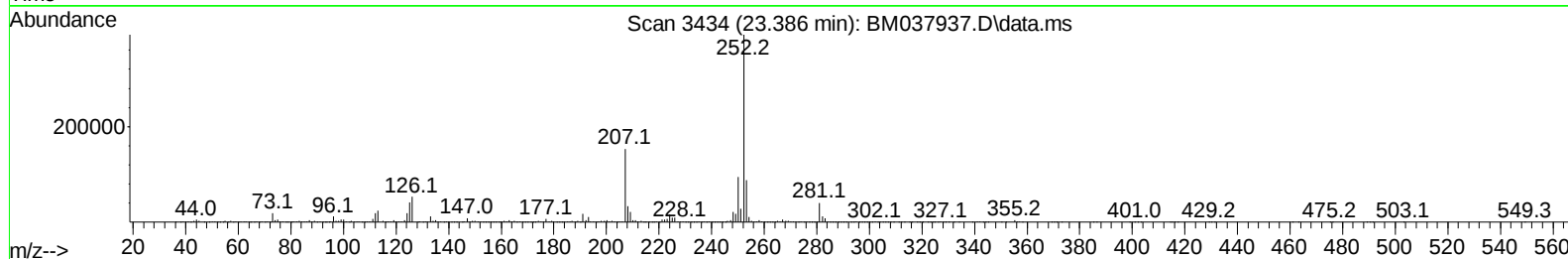
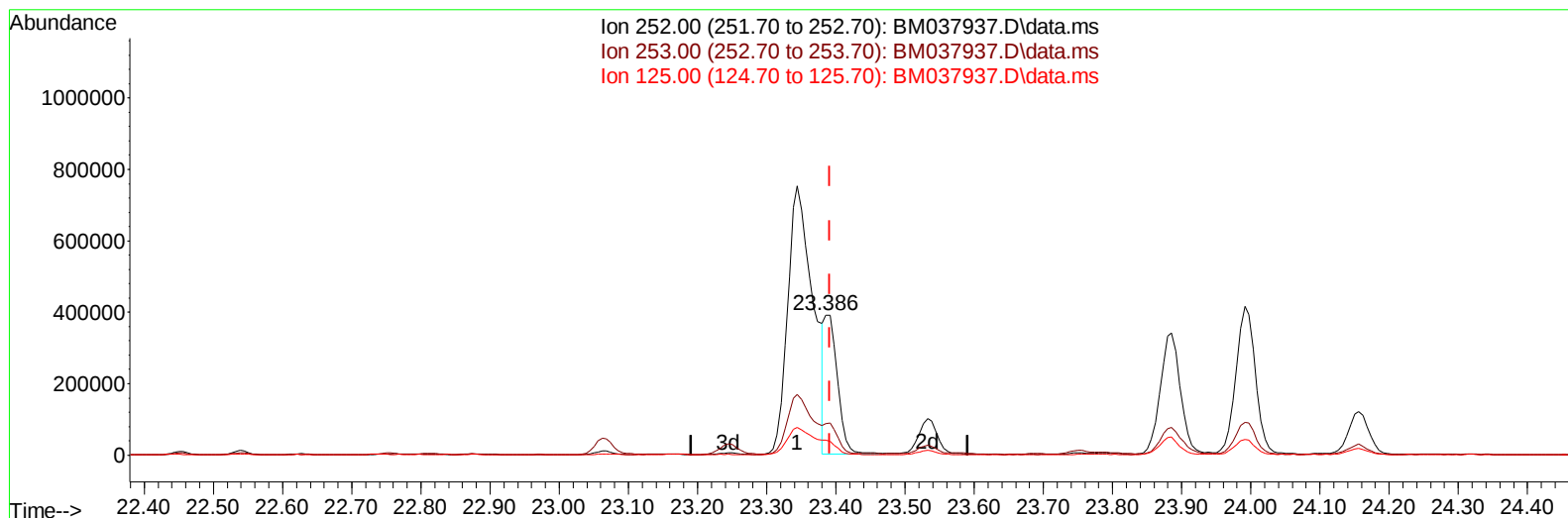
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(91) Benzo(k)fluoranthene

23.386min (-0.006) 8.60 ng/ul m

response 510634

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.80	22.32
125.00	10.10	10.49
0.00	0.00	0.00

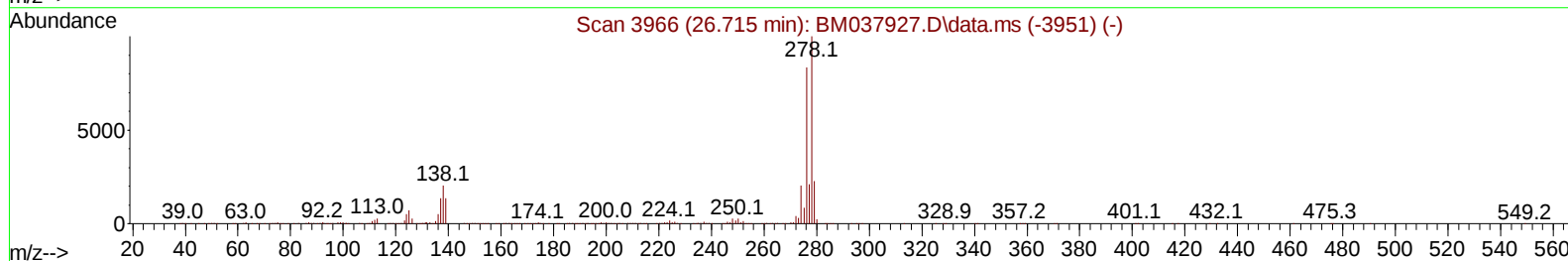
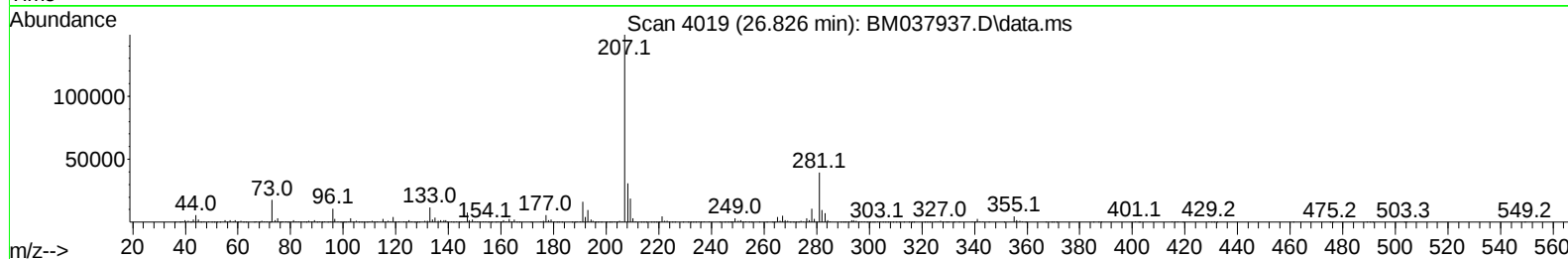
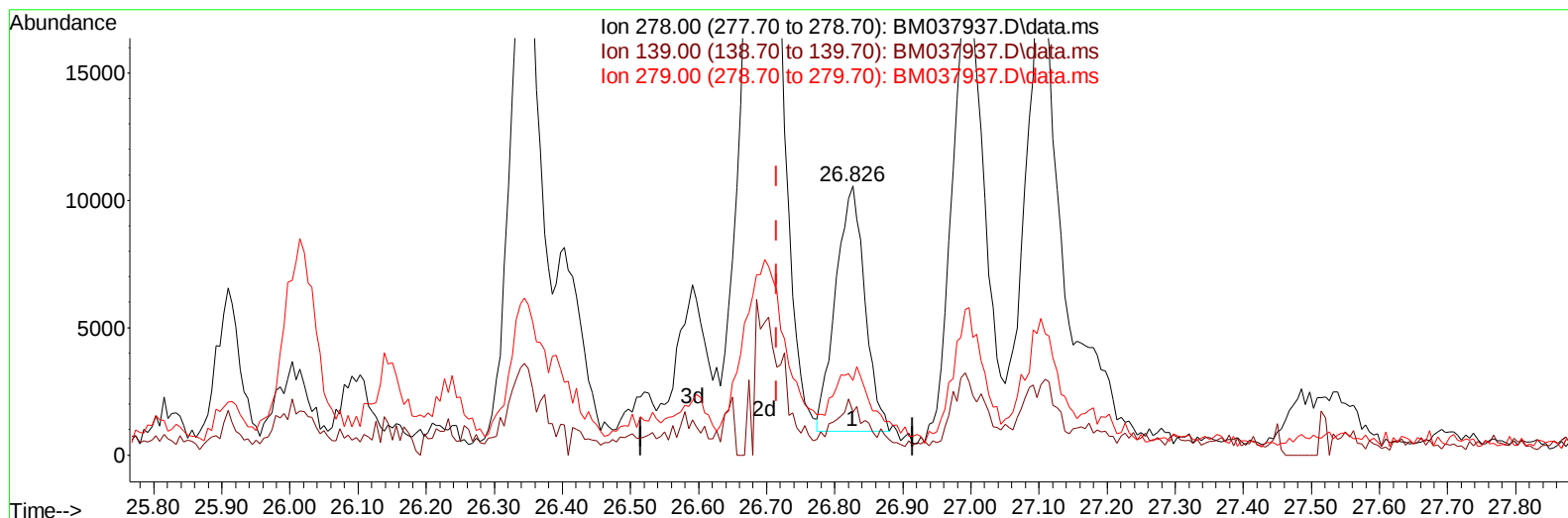
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(95) Dibenzo(a,h)anthracene

26.826min (+ 0.112) 0.49 ng/u1

response 28434

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	15.50	15.94
279.00	24.70	27.62
0.00	0.00	0.00

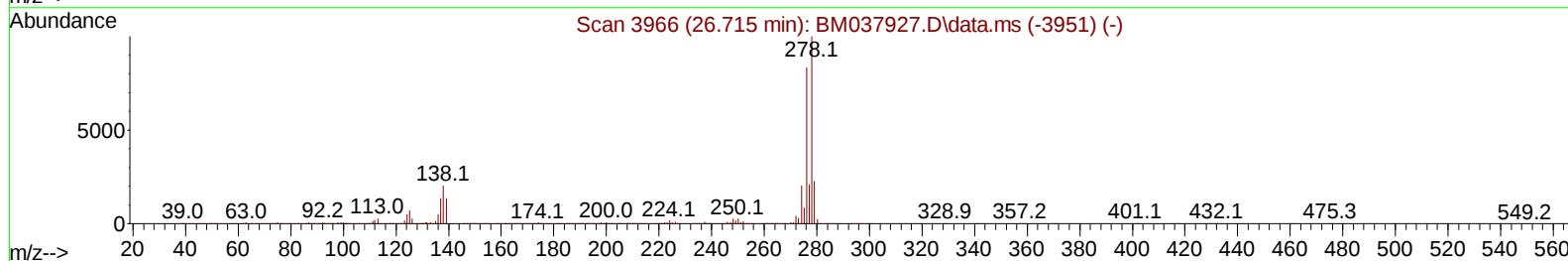
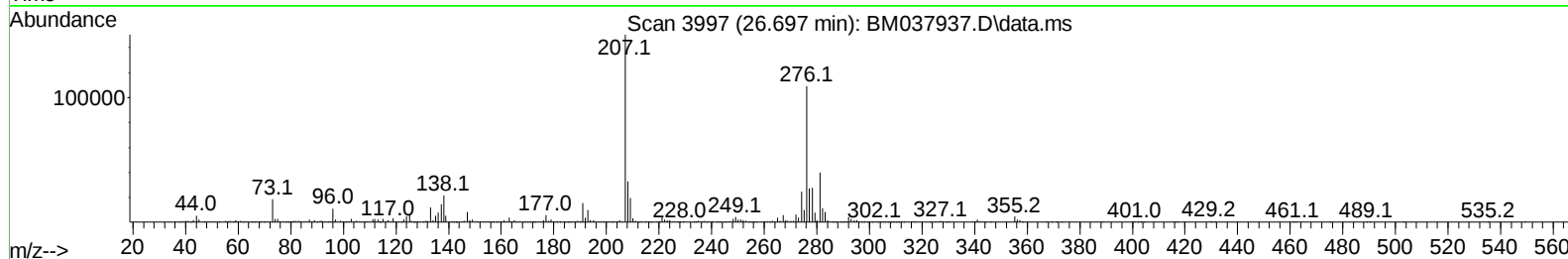
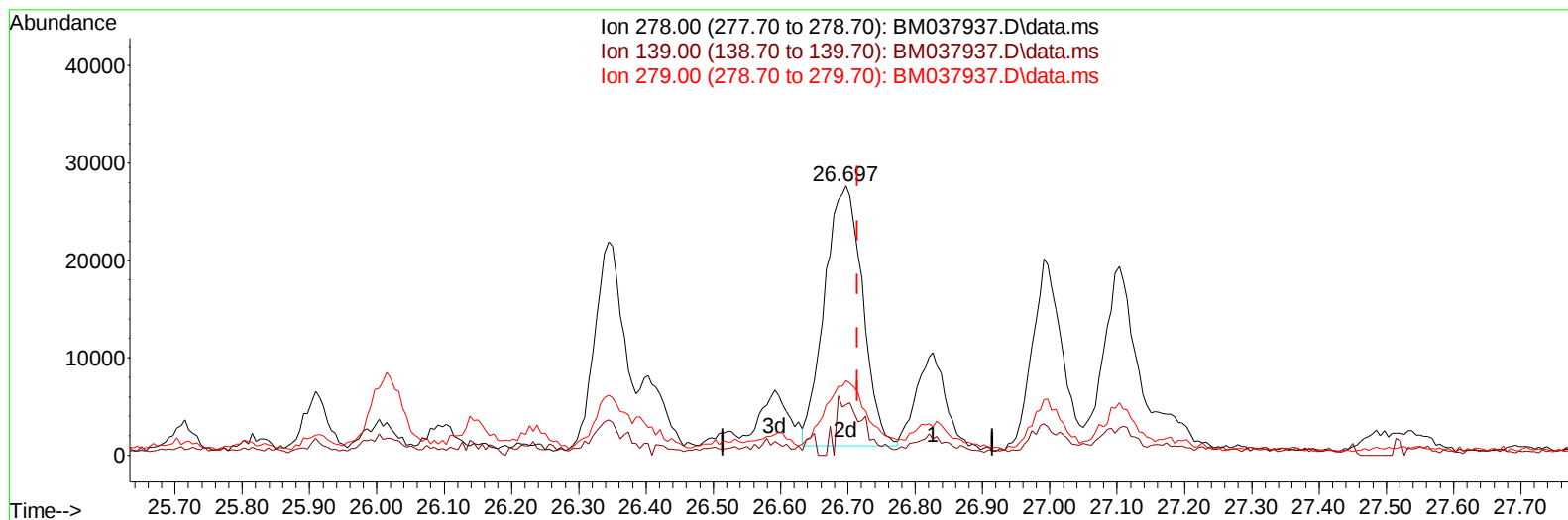
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(95) Dibenzo(a,h)anthracene

26.697min (-0.018) 1.78 ng/ul m

response 104104

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	15.50	18.86#
279.00	24.70	27.78
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.075	152	227991	20.000	ng/ul	0.00
20) Naphthalene-d8	10.898	136	848207	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.710	164	468407	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.451	188	925036	20.000	ng/ul	0.00
79) Chrysene-d12	21.621	240	870930	20.000	ng/ul	0.00
88) Perylene-d12	24.103	264	988297	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.422	96	24282	4.836	ng/uL	0.00
4) Pyridine-d5	3.857	84	207921	13.959	ng/ul	0.00
7) Phenol-d5	7.234	99	480409	25.931	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.398	67	308764	27.370	ng/ul	0.00
11) 2-Chlorophenol-d4	7.604	132	428000	28.161	ng/ul	0.00
15) 4-Methylphenol-d8	8.786	113	350352	24.226	ng/ul	0.00
21) Nitrobenzene-d5	9.251	128	206549	30.756	ng/ul	0.00
24) 2-Nitrophenol-d4	9.975	143	219172	31.142	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.516	165	368568	27.572	ng/ul	0.00
31) 4-Chloroaniline-d4	11.033	131	432399	23.433	ng/ul	0.00
46) Dimethylphthalate-d6	14.116	166	1044905	31.290	ng/ul	0.00
49) Acenaphthylene-d8	14.404	160	1231050	30.162	ng/ul	0.00
54) 4-Nitrophenol-d4	14.898	143	124542	21.883	ng/ul	0.00
60) Fluorene-d10	15.698	176	931875	31.322	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.821	200	114354	23.122	ng/ul	0.00
73) Anthracene-d10	17.551	188	1340893	31.059	ng/ul	0.00
81) Pyrene-d10	19.833	212	1566649	29.945	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.944	264	1610743	32.630	ng/ul	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	10.951	128	64710	1.414	ng/ul#	96
36) 2-Methylnaphthalene	12.545	142	42986	1.448	ng/ul	99
37) 1-Methylnaphthalene	12.763	142	211769	6.976	ng/ul	99
50) Acenaphthylene	14.433	152	140152	3.123	ng/ul#	95
52) Acenaphthene	14.774	153	2254853	72.641	ng/ul	100
61) Fluorene	15.757	166	2673324	77.349	ng/ul	98
72) Phenanthrene	17.498	178	9603986	190.117	ng/ul	97
74) Anthracene	17.592	178	8103437	157.847	ng/ul	99
80) Fluoranthene	19.509	202	9383568	151.976	ng/ul	99
82) Pyrene	19.868	202	7079410	110.169	ng/ul	99
85) Benzo(a)anthracene	21.603	228	1954923	33.341	ng/ul	99
87) Chrysene	21.662	228	2157168	39.212	ng/ul	98
90) Benzo(b)fluoranthene	23.344	252	1900536	31.019	ng/ul	99
91) Benzo(k)fluoranthene	23.386	252	510634m	8.599	ng/ul	
93) Benzo(a)pyrene	23.991	252	810965	15.051	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	26.691	276	381142	5.493	ng/ul	95
95) Dibenzo(a,h)anthracene	26.697	278	104104m	1.780	ng/ul	
96) Benzo(g,h,i)perylene	27.497	276	254257	4.589	ng/ul	97

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

