

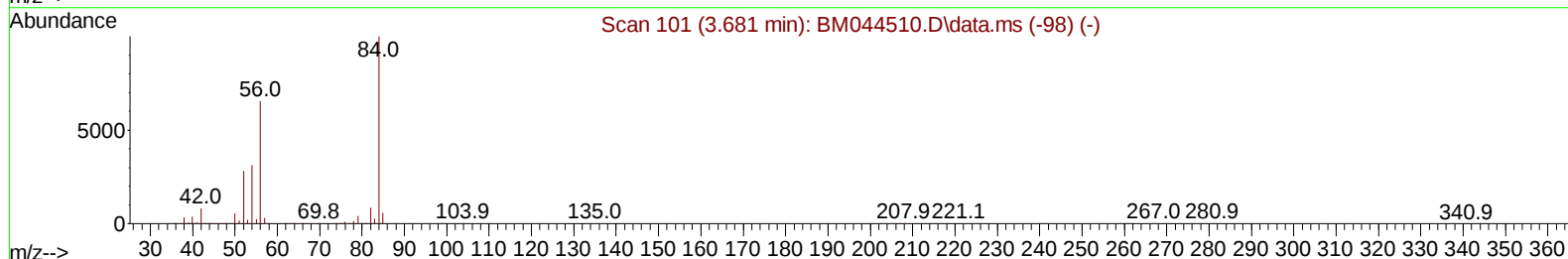
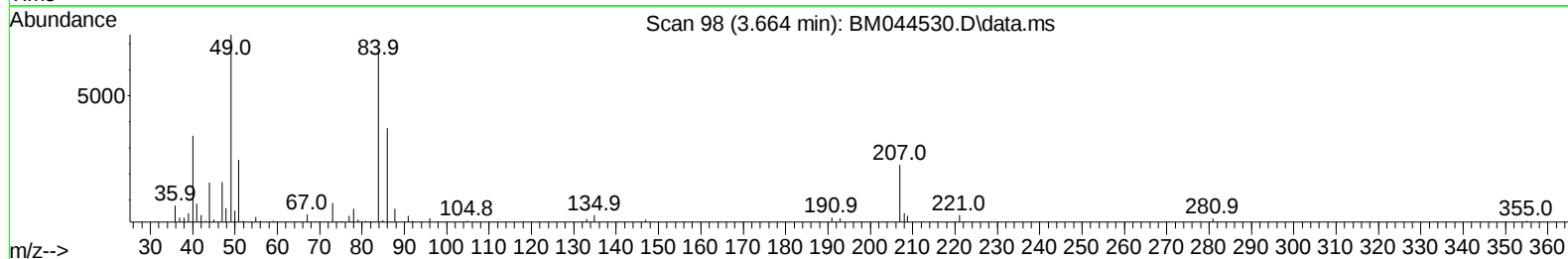
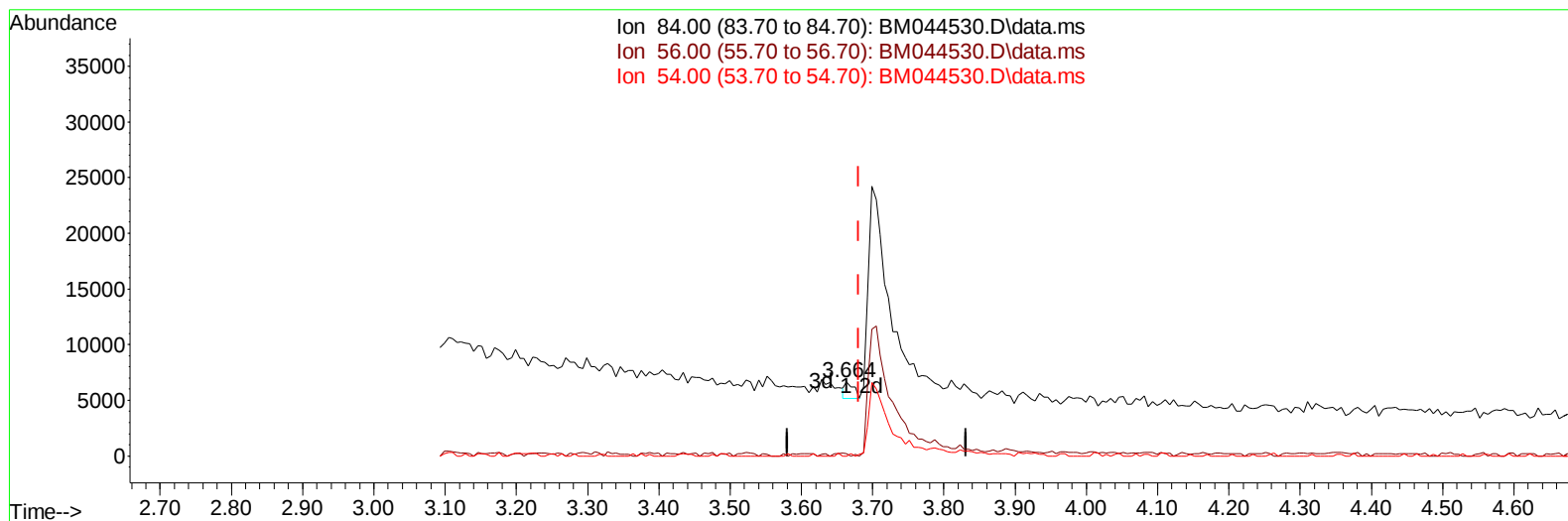
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
 Data File : BM044530.D
 Acq On : 06 Mar 2024 23:44
 Operator : MA/JU
 Sample : P1676-07
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E06Y6

Manual IntegrationsAPPROVED

Quant Time: Mar 07 00:19:42 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 05 23:05:11 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/07/2024
 Supervised By :mohammad ahmed 03/09/2024



TIC: BM044530.D\data.ms

(4) Pyridine-d5 (S)

3.664min (-0.017) 0.06 ng/u1

response 1237

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.10	3.21#
54.00	29.60	0.00#
0.00	0.00	0.00

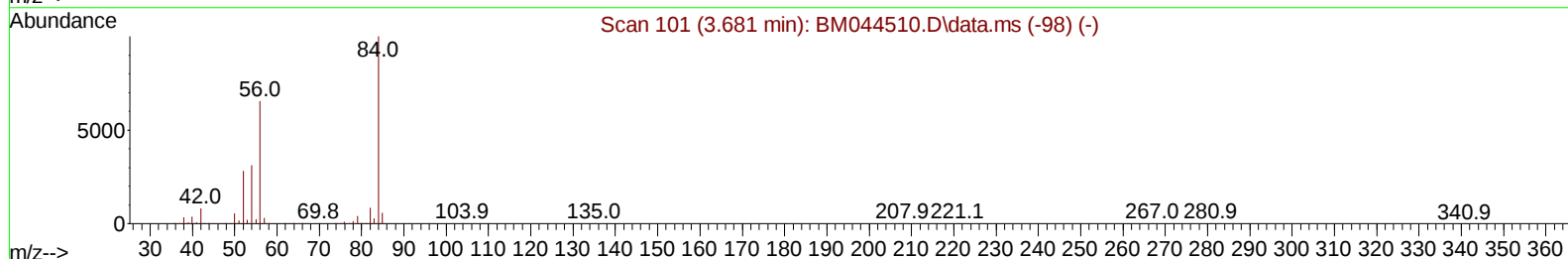
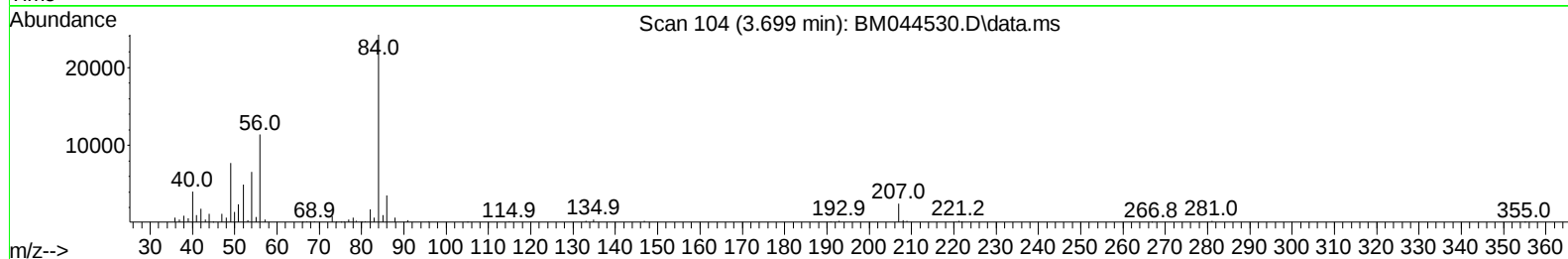
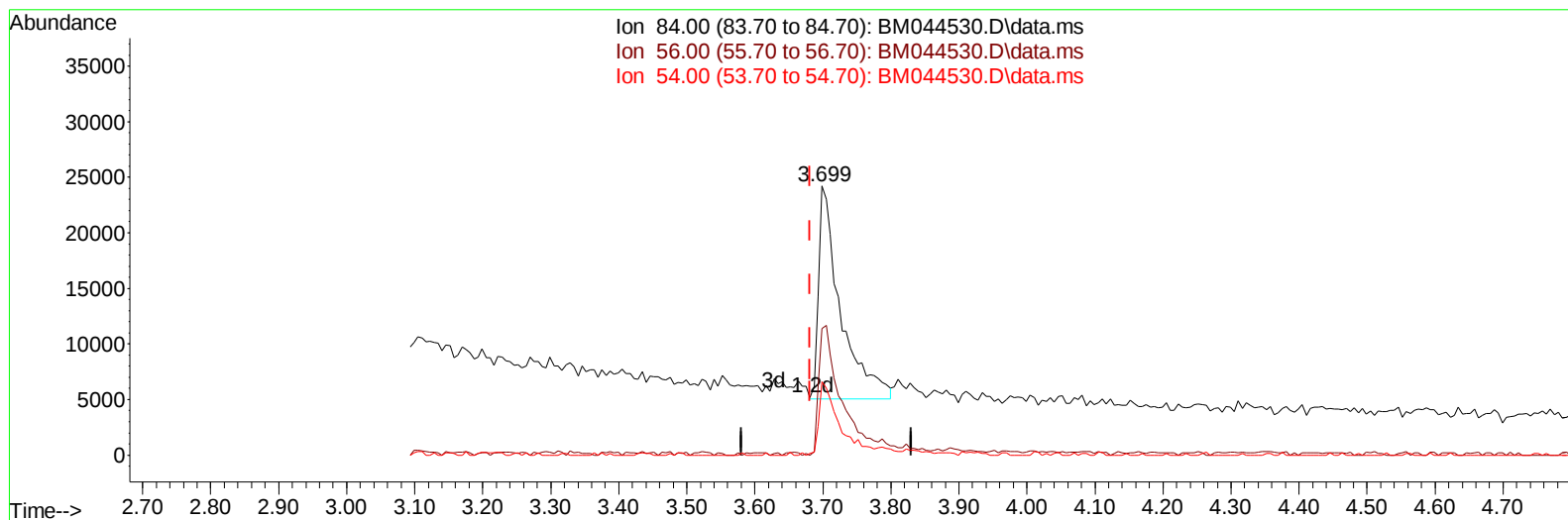
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(4) Pyridine-d5 (S)

3.699min (+ 0.018) 2.08 ng/u1 m

response 42528

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.10	46.98#
54.00	29.60	27.36
0.00	0.00	0.00

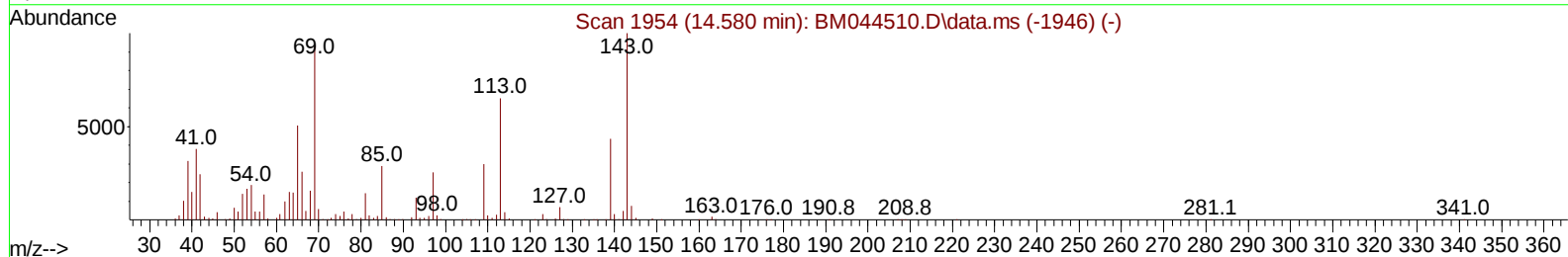
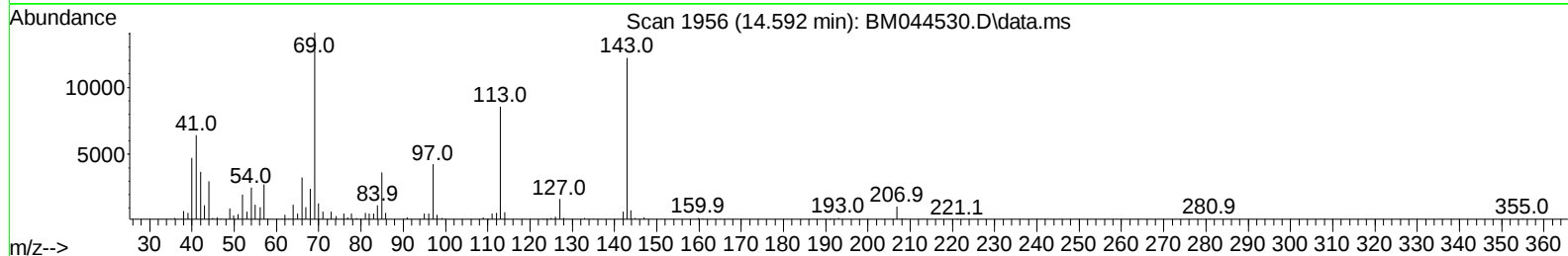
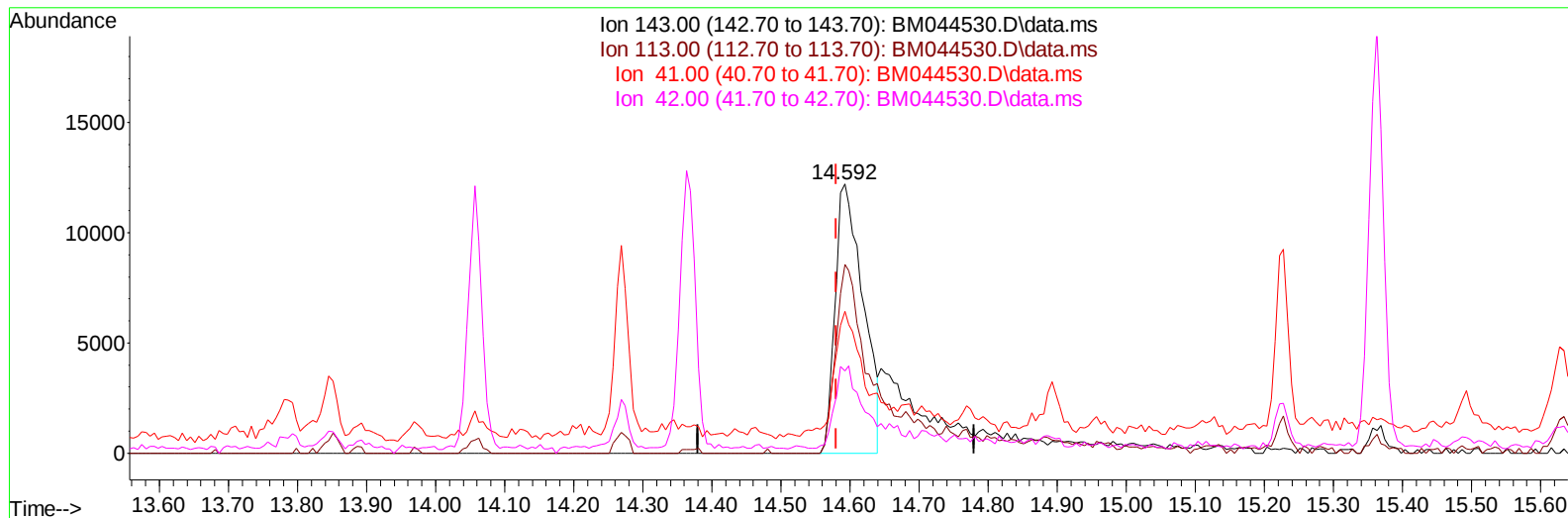
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Sample : P1676-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E06Y6

Manual IntegrationsAPPROVED

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TIC: BM044530.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.592min (+ 0.012) 3.14 ng/u1

response 34139

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	63.70	70.20
41.00	38.40	52.81#
42.00	26.60	30.45

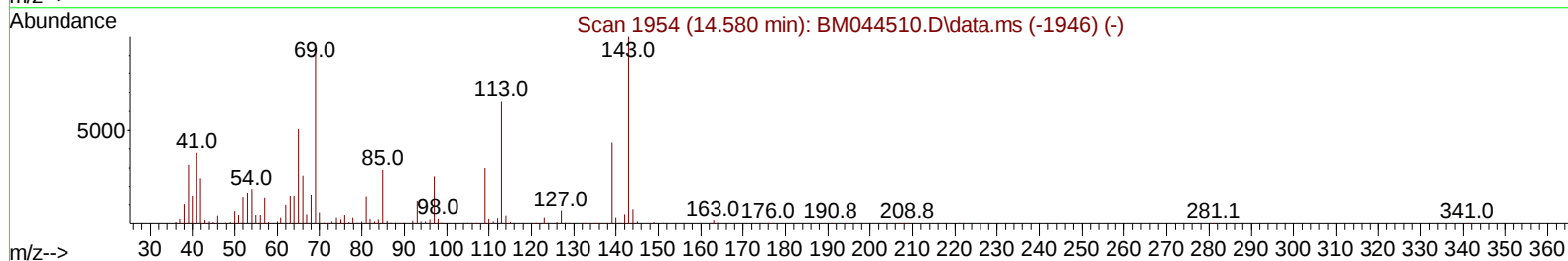
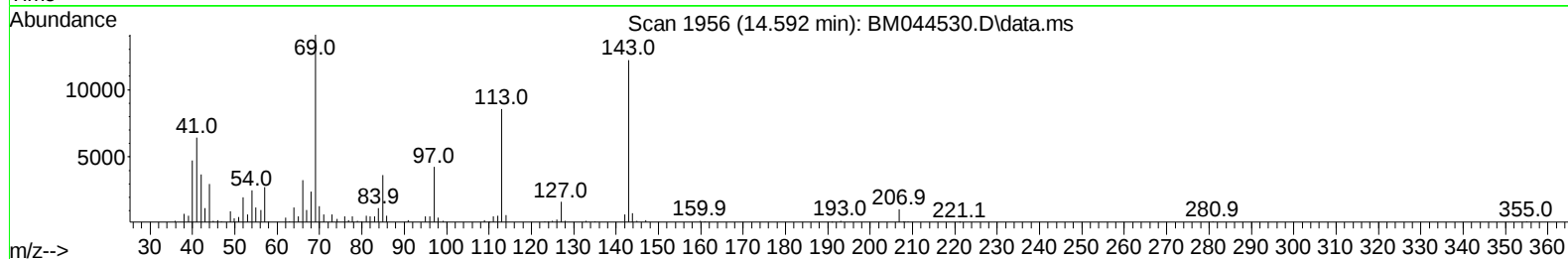
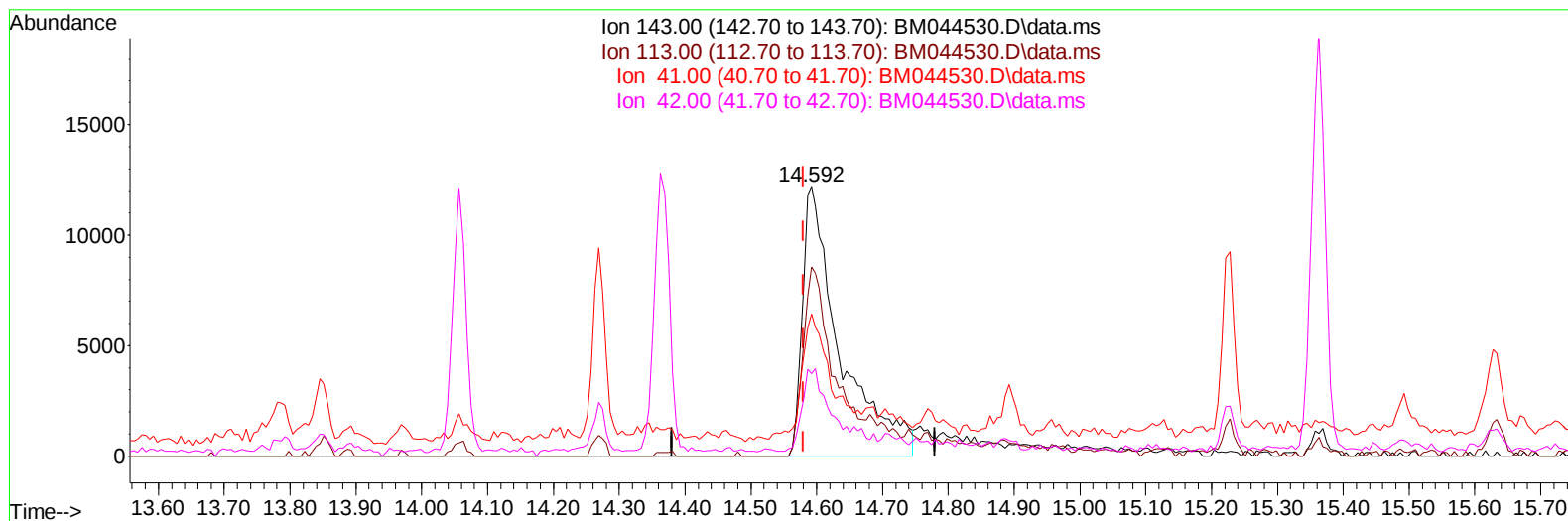
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Operator : MA/JU
Sample : P1676-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BM044530.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.592min (+ 0.012) 4.45 ng/ul m

response 48318

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	63.70	70.20
41.00	38.40	52.81#
42.00	26.60	30.45

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	236093	20.000	ng/ul	0.00
20) Naphthalene-d8	10.510	136	1034555	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.363	164	667954	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.116	188	1355041	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	1076461	20.000	ng/ul	0.00
88) Perylene-d12	23.551	264	1047215	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	16171	2.323	ng/uL	0.00
4) Pyridine-d5	3.699	84	42528m	2.076	ng/ul	0.02
7) Phenol-d5	6.899	99	378787	15.605	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.081	67	294469	19.640	ng/ul	0.00
11) 2-Chlorophenol-d4	7.263	132	294886	16.532	ng/ul	0.00
15) 4-Methylphenol-d8	8.434	113	310887	16.309	ng/ul	0.00
21) Nitrobenzene-d5	8.887	128	180470	20.716	ng/ul	0.00
24) 2-Nitrophenol-d4	9.604	143	118800	12.650	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	272983	17.079	ng/ul	0.00
31) 4-Chloroaniline-d4	10.657	131	331182	12.476	ng/ul	0.00
46) Dimethylphthalate-d6	13.786	166	838490	16.003	ng/ul	0.00
49) Acenaphthylene-d8	14.057	160	1404789	23.443	ng/ul	0.00
54) 4-Nitrophenol-d4	14.592	143	48318m	4.447	ng/ul	0.01
60) Fluorene-d10	15.363	176	1109606	25.200	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	21044	2.392	ng/ul	0.00
73) Anthracene-d10	17.216	188	1691335	26.113	ng/ul	0.00
81) Pyrene-d10	19.510	212	1915474	29.012	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.409	264	1446526	26.280	ng/ul	-0.01

Target Compounds Qvalue

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

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