

Quantitation Report (Qedit)

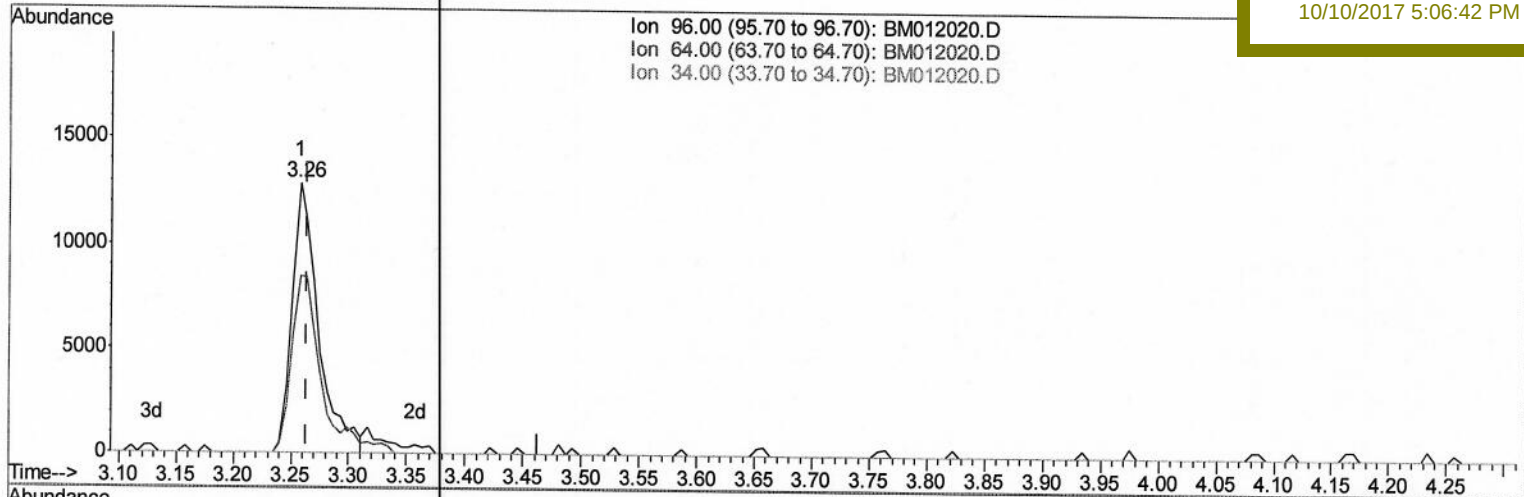
Data Path : Z:\HPCHEM1\BNA M\DATA\BM100917\
 Data File : BM012020.D
 Acq On : 09 Oct 2017 19:37
 Operator : SJ/JU
 Sample : I5666-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Oct 10 01:18:35 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 10 01:14:20 2017
 Response via : Initial Calibration

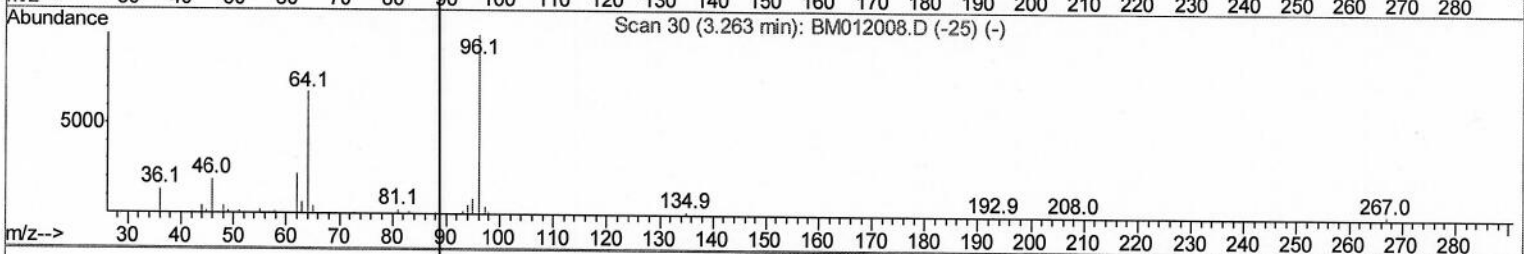
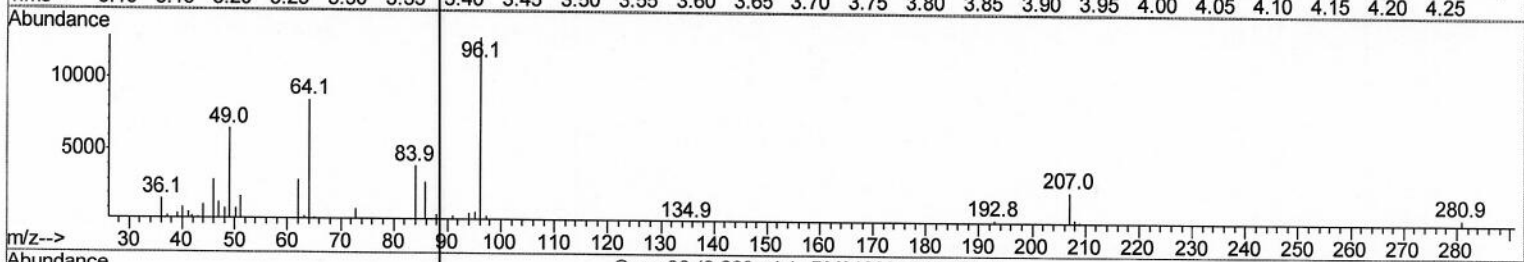
Instrument :
 BNA_M
 ClientSampleId :
 H0AD5

Manual Integrations
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Ion 96.00 (95.70 to 96.70): BM012020.D
 Ion 64.00 (63.70 to 64.70): BM012020.D
 Ion 34.00 (33.70 to 34.70): BM012020.D



TIC: BM012020.D

(3) 1,4-Dioxane-d8 (S)
 3.258min (-0.006) 7.30ng/uL
 response 20623

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	66.85
34.00	0.00	0.00
0.00	0.00	0.00

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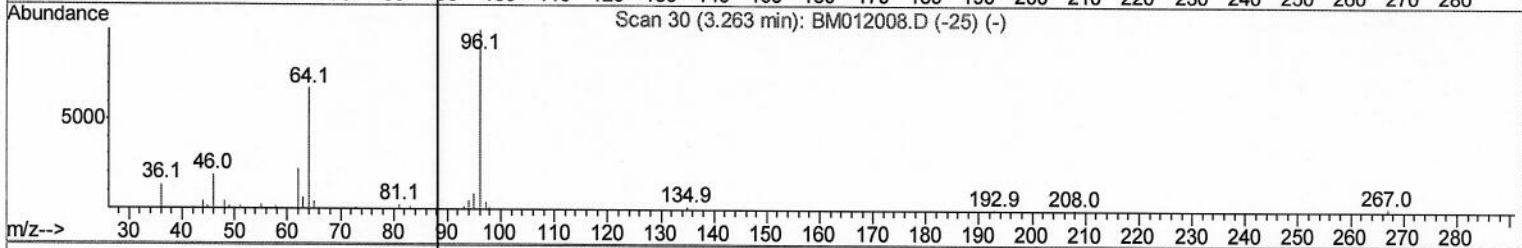
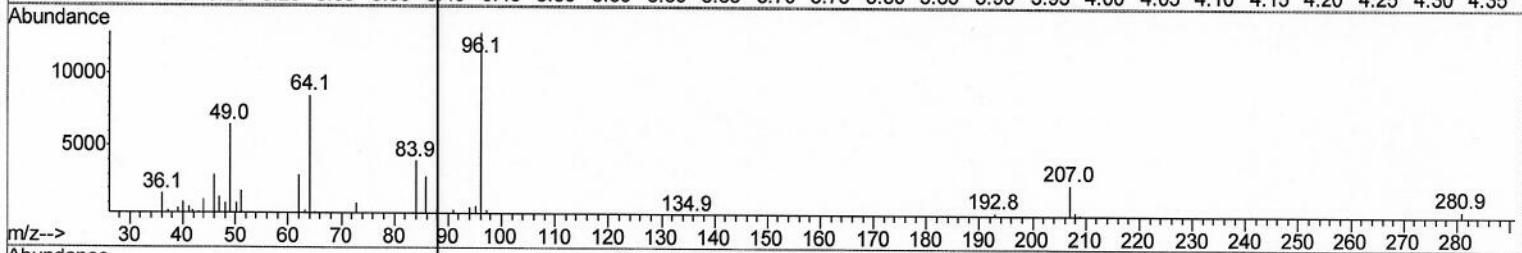
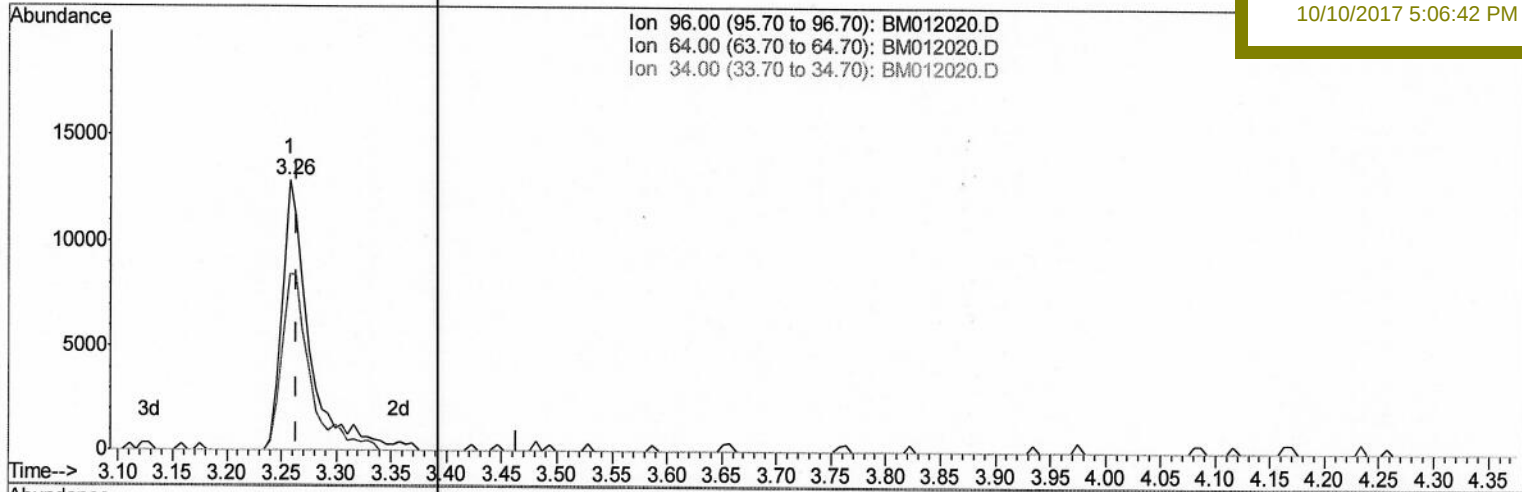
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TIC: BM012020.D

(3) 1,4-Dioxane-d8 (S)

3.258min (-0.006) 7.95ng/uL m

response 22476

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	61.34
34.00	0.00	0.00
0.00	0.00	0.00

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Quantitation Report (Qedit)

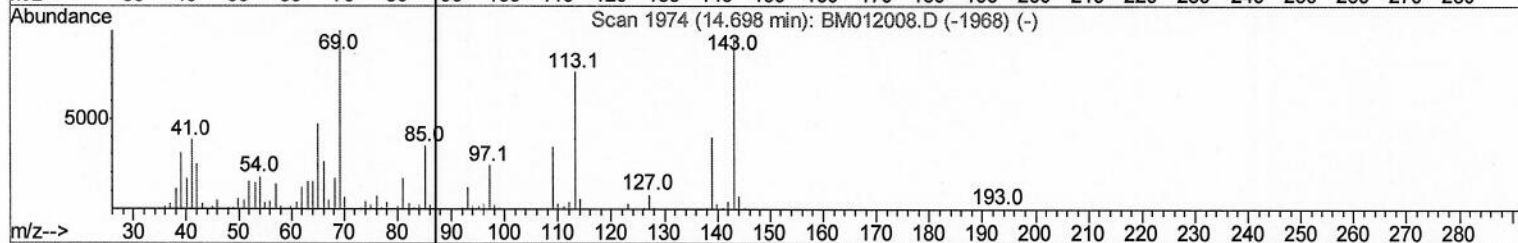
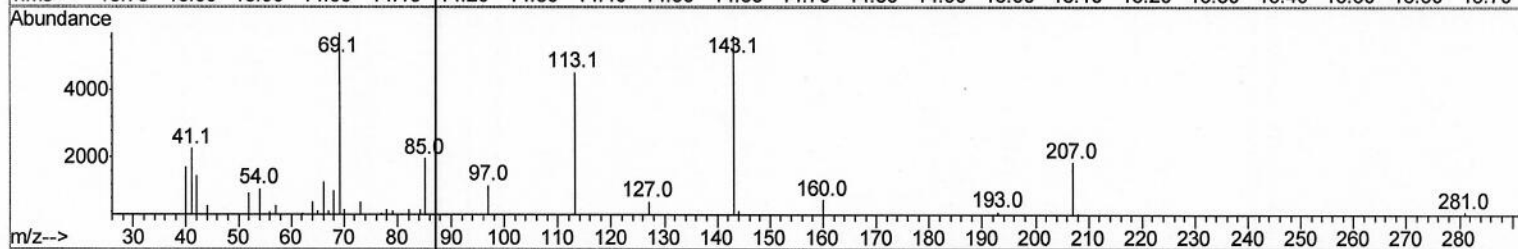
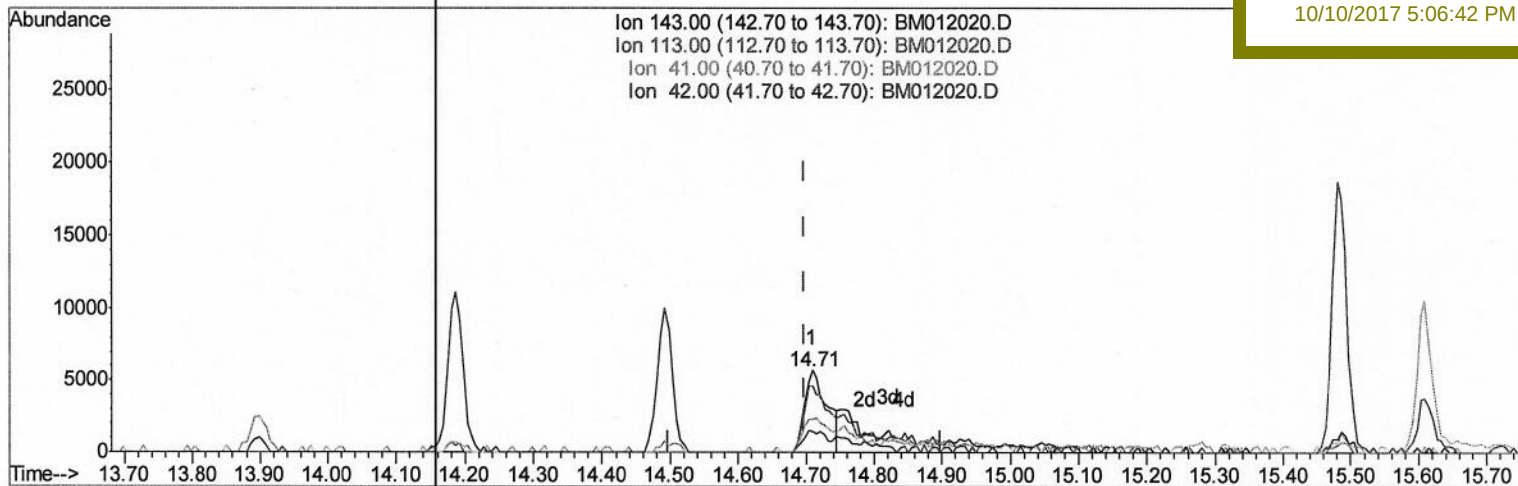
Data Path : Z:\HPCHEM1\BNA M\DATA\BM100917\
 Data File : BM012020.D
 Acq On : 09 Oct 2017 19:37
 Operator : SJ/JU
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 Misc :
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Quant Time: Oct 10 02:11:06 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 10 01:14:20 2017
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Instrument :
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TIC: BM012020.D

(51) 4-Nitrophenol-d4 (S)
 14.710min (+0.012) 2.16ng/ul
 response 13188

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	80.39#
41.00	34.80	40.06
42.00	20.50	26.03#

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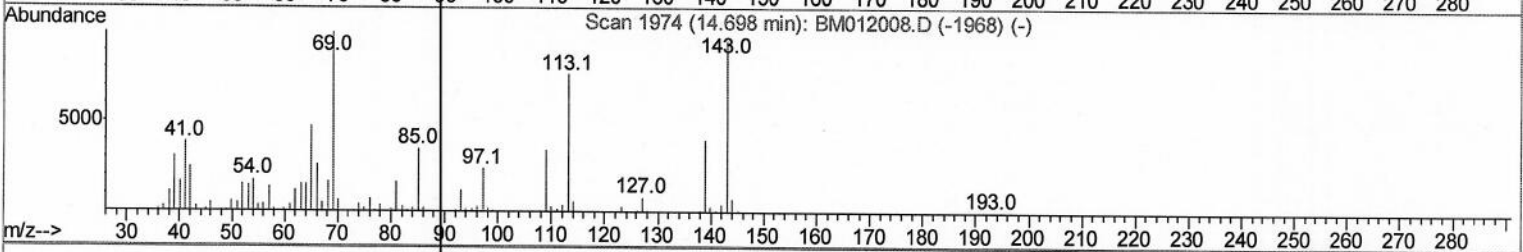
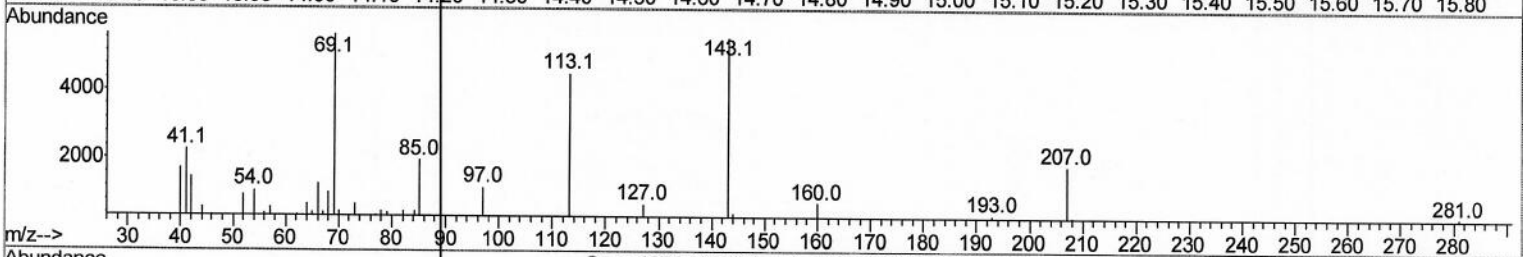
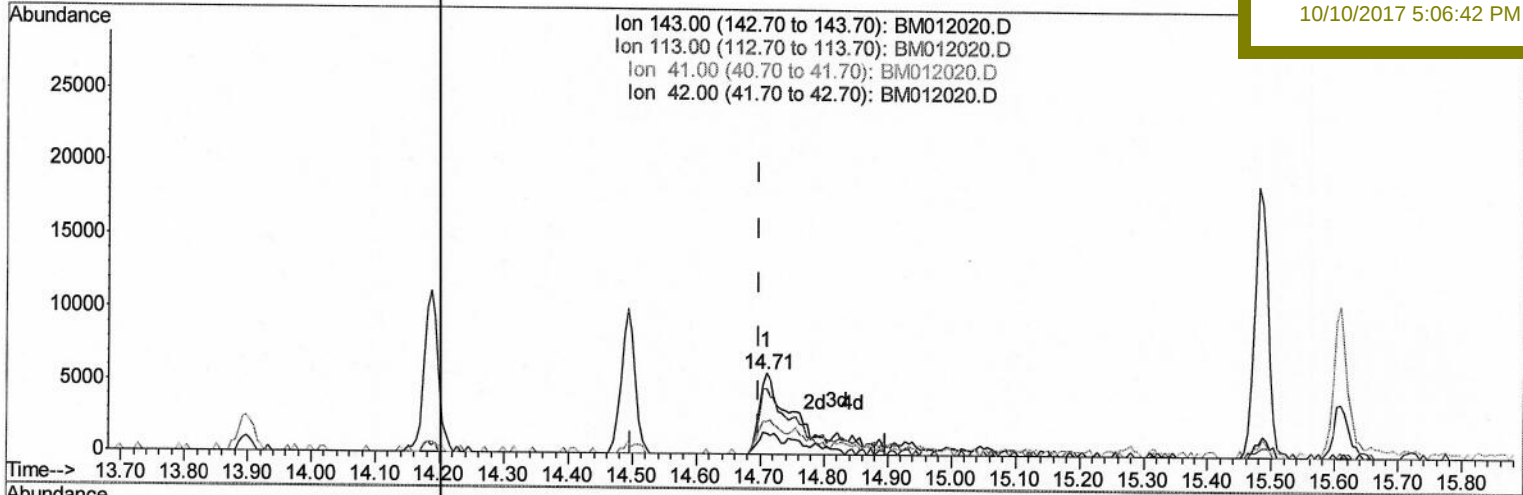
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TIC: BM012020.D

(51) 4-Nitrophenol-d4 (S)

14.710min (+0.012) 4.00ng/ul m

response 24409

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	80.39#
41.00	34.80	40.06
42.00	20.50	26.03#

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	133650	20.00	nq/ul	0.00
18) Naphthalene-d8	10.65	136	611557	20.00	nq/ul	0.00
35) Acenaphthene-d10	14.49	164	402808	20.00	nq/ul	0.00
61) Phenanthrene-d10	17.24	188	1009334	20.00	nq/ul	0.00
75) Chrysene-d12	21.41	240	1388202	20.00	nq/ul	0.00
83) Perylene-d12	23.73	264	1682645	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.26	96	22476m	7.95	nq/uL	0.00
5) Phenol-d5	7.01	99	77577	6.75	nq/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.17	67	197036	28.95	nq/ul	0.00
9) 2-Chlorophenol-d4	7.37	132	229592	24.89	nq/ul	0.00
13) 4-Methylphenol-d8	8.55	113	154607	15.55	nq/ul	0.00
19) Nitrobenzene-d5	9.02	128	132956	30.53	nq/ul	0.00
22) 2-Nitrophenol-d4	9.73	143	150185	30.62	nq/ul	0.00
26) 2,4-Dichlorophenol-d3	10.27	165	282218	28.47	nq/ul	0.00
29) 4-Chloroaniline-d4	10.82	131	3309	0.31	nq/ul	0.02
43) Dimethylphthalate-d6	13.90	166	1163686	33.33	nq/ul	0.00
46) Acenaphthylene-d8	14.19	160	1346643	32.83	nq/ul	0.00
51) 4-Nitrophenol-d4	14.71	143	24409m	4.00	nq/ul	0.01
57) Fluorene-d10	15.49	176	999397	32.44	nq/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.61	200	192032	28.28	nq/ul	0.00
70) Anthracene-d10	17.33	188	1727926	34.74	nq/ul	0.00
76) Pyrene-d10	19.63	212	2207878	35.26	nq/ul	0.00
87) Benzo(a)pyrene-d12	23.58	264	2789274	34.46	ng/ul	0.00

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Target Compounds

Ovalue

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

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