

Data Path : Z:\HPCHEM1\BNA M\DATA\BM043016\
 Data File : BM005148.D
 Acq On : 29 Apr 2016 18:31
 Operator : UM/SJ
 Sample : H2778-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B68

Quant Time: Apr 30 04:05:47 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM042516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Apr 30 03:07:27 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	38012	20.00	ng/ul	0.00
18) Naphthalene-d8	10.57	136	181787	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.43	164	117878	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.17	188	277825	20.00	ng/ul	0.00
75) Chrysene-d12	21.36	240	304988	20.00	ng/ul	0.00
83) Perylene-d12	23.64	264	251314	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	788	1.05	ng/uL	0.00
5) Phenol-d5	6.96	99	15889	5.05	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	55143	30.82	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	55649	22.37	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	34099	12.79	ng/ul	0.00
19) Nitrobenzene-d5	8.95	128	35955	28.56	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	40754	29.01	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	68858	25.80	ng/ul	0.00
29) 4-Chloroaniline-d4	10.73	131	3191	1.02	ng/ul	0.01
43) Dimethylphthalate-d6	13.83	166	273947	29.43	ng/ul	0.00
46) Acenaphthylene-d8	14.12	160	325564	29.37	ng/ul	0.00
51) 4-Nitrophenol-d4	14.64	143	3692	2.27	ng/ul	0.00
57) Fluorene-d10	15.42	176	240354	29.38	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.55	200	33846	21.42	ng/ul	0.00
70) Anthracene-d10	17.27	188	376501	30.22	ng/ul	0.00
76) Pyrene-d10	19.57	212	425741	30.66	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.50	264	342393	30.41	ng/ul	0.00

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

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