

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM043016\
 Data File : BM005149.D
 Acq On : 29 Apr 2016 19:07
 Operator : UM/SJ
 Sample : H2778-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B69

Quant Time: Apr 30 03:09:49 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	39445	20.00	ng/ul	0.00
18) Naphthalene-d8	10.57	136	186360	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.43	164	120451	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.17	188	286992	20.00	ng/ul	0.00
75) Chrysene-d12	21.36	240	313267	20.00	ng/ul	0.00
83) Perylene-d12	23.64	264	260458	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	894	1.15	ng/uL	0.00
5) Phenol-d5	6.96	99	17724	5.43	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	56080	30.21	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	57702	22.35	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	36603	13.23	ng/ul	0.00
19) Nitrobenzene-d5	8.95	128	36307	28.13	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	41624	28.91	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	68764	25.14	ng/ul	0.00
29) 4-Chloroaniline-d4	10.73	131	2327	0.73	ng/ul	0.02
43) Dimethylphthalate-d6	13.83	166	274354	28.84	ng/ul	0.00
46) Acenaphthylene-d8	14.12	160	327772	28.93	ng/ul	0.00
51) 4-Nitrophenol-d4	14.64	143	3904	2.35	ng/ul	0.00
57) Fluorene-d10	15.42	176	240548	28.78	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.54	200	33176	20.32	ng/ul	0.00
70) Anthracene-d10	17.27	188	377598	29.34	ng/ul	0.00
76) Pyrene-d10	19.57	212	425954	29.87	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.50	264	342502	29.35	ng/ul	0.00

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

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