

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080616\
 Data File : BM006947.D
 Acq On : 06 Aug 2016 20:45
 Operator : UM/SJ
 Sample : H4301-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA0E9

Quant Time: Aug 08 03:53:00 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 08 03:51:46 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	130664	20.00	ng/ul	0.00
7) Naphthalene-d8	10.33	136	631247	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.21	164	433323	20.00	ng/ul	0.00
23) Phenanthrene-d10	16.96	188	1074724	20.00	ng/ul	0.00
29) Chrysene-d12	21.17	240	1391368	20.00	ng/ul	0.00
34) Perylene-d12	23.36	264	1309402	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.08	96	4152	1.45	ng/uL	0.00
3) Phenol-d5	6.78	99	190020	15.34	ng/ul	0.01
4) Bis-(2-Chloroethyl)ether-d	6.90	67	136752	17.60	ng/ul	0.00
5) 2-Chlorophenol-d4	7.10	132	149950	16.95	ng/ul	0.00
6) 4-Methylphenol-d8	8.30	113	145706	14.09	ng/ul	0.00
8) Nitrobenzene-d5	8.72	128	77956	16.63	ng/ul	0.00
9) 2-Nitrophenol-d4	9.44	143	91818	16.35	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	9.99	165	175635	16.50	ng/ul	0.02
12) 4-Chloroaniline-d4	10.50	131	224057	17.63	ng/ul	0.01
16) Dimethylphthalate-d6	13.63	166	728049	19.09	ng/ul	0.00
17) Acenaphthylene-d8	13.90	160	819641	18.58	ng/ul	0.00
20) 4-Nitrophenol-d4	14.53	143	57648	10.42	ng/ul	0.04
21) Fluorene-d10	15.21	176	626565	19.33	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.37	200	54561	12.30	ng/ul	0.00
26) Anthracene-d10	17.06	188	1012778	19.51	ng/ul	0.00
30) Pyrene-d10	19.37	212	1266436	19.09	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.22	264	1242019	19.92	ng/ul	0.00

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

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