

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080616\
 Data File : BM006949.D
 Acq On : 06 Aug 2016 21:57
 Operator : UM/SJ
 Sample : H4301-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA0F7

Quant Time: Aug 08 03:53:16 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	158612	20.00	ng/ul	0.00
7) Naphthalene-d8	10.33	136	760753	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.20	164	530093	20.00	ng/ul	0.00
23) Phenanthrene-d10	16.96	188	1258707	20.00	ng/ul	0.00
29) Chrysene-d12	21.17	240	1490749	20.00	ng/ul	0.00
34) Perylene-d12	23.36	264	1256913	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.09	96	5665	1.63	ng/uL	0.00
3) Phenol-d5	6.77	99	223425	14.86	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.90	67	168773	17.90	ng/ul	0.00
5) 2-Chlorophenol-d4	7.10	132	191790	17.86	ng/ul	0.00
6) 4-Methylphenol-d8	8.30	113	160865	12.81	ng/ul	0.00
8) Nitrobenzene-d5	8.72	128	103673	18.35	ng/ul	0.00
9) 2-Nitrophenol-d4	9.43	143	122862	18.16	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	9.99	165	205733	16.04	ng/ul	0.01
12) 4-Chloroaniline-d4	10.50	131	287816	18.79	ng/ul	0.00
16) Dimethylphthalate-d6	13.63	166	904358	19.39	ng/ul	0.00
17) Acenaphthylene-d8	13.90	160	1029185	19.07	ng/ul	0.00
20) 4-Nitrophenol-d4	14.53	143	85962	12.70	ng/ul	0.04
21) Fluorene-d10	15.21	176	773084	19.50	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.37	200	69098	13.29	ng/ul	0.00
26) Anthracene-d10	17.06	188	1198968	19.72	ng/ul	0.00
30) Pyrene-d10	19.37	212	1438621	20.24	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.22	264	1188612	19.85	ng/ul	0.00

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

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