

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072016\
 Data File : BM006572.D
 Acq On : 21 Jul 2016 01:50
 Operator : UM/SJ
 Sample : H4076-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9YP6

Quant Time: Jul 21 03:39:47 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 21 01:09:30 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	184245	20.00	ng/ul	0.00
7) Naphthalene-d8	10.39	136	795366	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.26	164	463494	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.02	188	1051601	20.00	ng/ul	0.00
29) Chrysene-d12	21.21	240	1227106	20.00	ng/ul	0.00
34) Perylene-d12	23.40	264	1278562	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.14	96	8894	1.95	ng/uL	0.00
3) Phenol-d5	6.79	99	343143	21.89	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.95	67	220878	22.75	ng/ul	0.00
5) 2-Chlorophenol-d4	7.15	132	262523	20.96	ng/ul	0.00
6) 4-Methylphenol-d8	8.32	113	266965	21.96	ng/ul	0.00
8) Nitrobenzene-d5	8.77	128	128740	21.38	ng/ul	0.00
9) 2-Nitrophenol-d4	9.49	143	137347	21.45	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.02	165	253508	20.95	ng/ul	0.00
12) 4-Chloroaniline-d4	10.53	131	335693	25.19	ng/ul	0.00
16) Dimethylphthalate-d6	13.67	166	769870	21.18	ng/ul	0.00
17) Acenaphthylene-d8	13.95	160	982952	21.19	ng/ul	0.00
20) 4-Nitrophenol-d4	14.48	143	111664	17.69	ng/ul	0.00
21) Fluorene-d10	15.26	176	688602	21.19	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.39	200	89778	16.45	ng/ul	0.00
26) Anthracene-d10	17.11	188	1070814	21.45	ng/ul	0.00
30) Pyrene-d10	19.42	212	1237021	22.12	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.27	264	1303617	22.17	ng/ul	0.00

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

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