

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072016\
 Data File : BM006573.D
 Acq On : 21 Jul 2016 02:27
 Operator : UM/SJ
 Sample : H4078-16
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9YW0

Quant Time: Jul 21 03:39:53 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	7.61	152	180110	20.00	ng/ul	0.00
7) Naphthalene-d8	10.39	136	751278	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.26	164	414872	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.02	188	942690	20.00	ng/ul	0.00
29) Chrysene-d12	21.21	240	1122323	20.00	ng/ul	0.00
34) Perylene-d12	23.41	264	1197254	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.14	96	2234	0.50	ng/uL	0.00
3) Phenol-d5	6.79	99	76267	4.98	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.95	67	221889	23.38	ng/ul	0.00
5) 2-Chlorophenol-d4	7.15	132	227115	18.55	ng/ul	0.00
6) 4-Methylphenol-d8	8.32	113	139288	11.72	ng/ul	0.00
8) Nitrobenzene-d5	8.77	128	132669	23.33	ng/ul	0.00
9) 2-Nitrophenol-d4	9.48	143	143582	23.74	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.02	165	235208	20.58	ng/ul	0.00
12) 4-Chloroaniline-d4	10.55	131	11103	0.88	ng/ul	0.01
16) Dimethylphthalate-d6	13.67	166	833353	25.62	ng/ul	0.00
17) Acenaphthylene-d8	13.95	160	1022414	24.62	ng/ul	0.00
20) 4-Nitrophenol-d4	14.49	143	20886	3.70	ng/ul	0.00
21) Fluorene-d10	15.26	176	749408	25.76	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.39	200	111681	22.82	ng/ul	0.00
26) Anthracene-d10	17.11	188	1270867	28.39	ng/ul	0.00
30) Pyrene-d10	19.42	212	1426661	27.90	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.27	264	1587857	28.83	ng/ul	0.00

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

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