

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM071616\
 Data File : BM006493.D
 Acq On : 17 Jul 2016 00:17
 Operator : UM/SJ
 Sample : H3985-07
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BDJ46

Quant Time: Jul 18 07:26:53 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM071416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 15 23:20:22 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	184154	20.00	ng/ul	0.00
18) Naphthalene-d8	10.39	136	771839	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.27	164	428322	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.02	188	961668	20.00	ng/ul	0.00
75) Chrysene-d12	21.22	240	1170879	20.00	ng/ul	-0.01
83) Perylene-d12	23.41	264	1253631	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.14	96	12139	2.78	ng/uL	0.00
5) Phenol-d5	6.79	99	543647	36.02	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.96	67	336640	36.78	ng/ul	0.00
9) 2-Chlorophenol-d4	7.15	132	422080	34.52	ng/ul	0.00
13) 4-Methylphenol-d8	8.32	113	429186	35.79	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	208877	35.71	ng/ul	0.00
22) 2-Nitrophenol-d4	9.49	143	232332	35.13	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.02	165	405144	32.47	ng/ul	0.00
29) 4-Chloroaniline-d4	10.53	131	544957	41.78	ng/ul	0.00
43) Dimethylphthalate-d6	13.68	166	1263958	36.06	ng/ul	0.00
46) Acenaphthylene-d8	13.96	160	1573977	36.32	ng/ul	0.00
51) 4-Nitrophenol-d4	14.49	143	228312	37.49	ng/ul	0.00
57) Fluorene-d10	15.26	176	1084409	34.80	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.40	200	179859	32.33	ng/ul	0.00
70) Anthracene-d10	17.12	188	1697696	37.35	ng/ul	0.00
76) Pyrene-d10	19.42	212	1942468	36.59	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.27	264	2174364	37.93	ng/ul	-0.01

Target Compounds						Qvalue
6) Phenol	6.82	94	16122	1.01	ng/ul	96
44) Dimethylphthalate	13.73	163	976832	27.47	ng/ul	99

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

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