

Data Path : Z:\HPCHEM1\BNA M\DATA\BM080416\
 Data File : BM006870.D
 Acq On : 04 Aug 2016 19:08
 Operator : UM/SJ
 Sample : H4095-11
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BDJG4

Manual Integrations
 APPROVED

umangi
 8/5/2016 7:00:24 PM

Quant Time: Aug 05 04:38:13 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM080316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 05 02:18:21 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	224365	20.00	ng/ul	0.00
18) Naphthalene-d8	10.34	136	953490	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.22	164	616167	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.98	188	1444884	20.00	ng/ul	0.00
75) Chrysene-d12	21.19	240	1655959	20.00	ng/ul	0.00
83) Perylene-d12	23.37	264	1423408	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.11	96	8506	2.08	ng/uL	0.00
5) Phenol-d5	6.79	99	393173	23.81	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.92	67	276999	25.11	ng/ul	0.00
9) 2-Chlorophenol-d4	7.12	132	331223	25.38	ng/ul	0.00
13) 4-Methylphenol-d8	8.32	113	315479	23.49	ng/ul	0.00
19) Nitrobenzene-d5	8.73	128	153314	23.52	ng/ul	0.00
22) 2-Nitrophenol-d4	9.45	143	178085	22.94	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.00	165	324905	20.25	ng/ul	0.00
29) 4-Chloroaniline-d4	10.52	131	401394	24.21	ng/ul	0.00
43) Dimethylphthalate-d6	13.65	166	1292516	25.47	ng/ul	0.00
46) Acenaphthylene-d8	13.92	160	1509006	24.34	ng/ul	0.00
51) 4-Nitrophenol-d4	14.53	143	147910m	13.29	ng/ul	0.00
57) Fluorene-d10	15.22	176	1143681	24.28	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.38	200	173278	18.37	ng/ul	0.00
70) Anthracene-d10	17.07	188	1782147m	25.68	ng/ul	0.00
76) Pyrene-d10	19.39	212	2159421	25.49	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.23	264	1825293	26.33	ng/ul	0.00

Target Compounds					Ovalue
44) Dimethylphthalate	13.69	163	694265	13.76	ng/ul 98

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

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