

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121416\
 Data File : BM008314.D
 Acq On : 14 Dec 2016 18:23
 Operator : UM/SJ
 Sample : H5996-07
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA866

Manual Integrations
 APPROVED

sohil
 12/15/2016 5:36:13 PM

Quant Time: Dec 15 02:35:52 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 15 02:28:30 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	113208	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	438771	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.31	164	293542	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.06	188	601618	20.00	ng/ul	0.00
75) Chrysene-d12	21.27	240	777138	20.00	ng/ul	0.00
83) Perylene-d12	23.49	264	798678	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.24	96	3118	1.30	ng/uL	0.00
5) Phenol-d5	6.87	99	39064	4.45	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.02	67	150523	29.82	ng/ul	0.00
9) 2-Chlorophenol-d4	7.21	132	136510	20.66	ng/ul	0.00
13) 4-Methylphenol-d8	8.40	113	50158	7.37	ng/ul	0.01
19) Nitrobenzene-d5	8.83	128	97935	30.89	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	106791	30.39	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.10	165	150037	23.30	ng/ul	0.02
29) 4-Chloroaniline-d4	10.70	131	2518m	0.34	ng/ul	0.10
43) Dimethylphthalate-d6	13.73	166	658980	33.77	ng/ul	0.00
46) Acenaphthylene-d8	14.00	160	752202	32.62	ng/ul	0.00
51) 4-Nitrophenol-d4	14.64	143	8717m	3.03	ng/ul	0.08
57) Fluorene-d10	15.31	176	556438	32.81	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.47	200	87309	24.38	ng/ul	0.00
70) Anthracene-d10	17.17	188	725985	27.70	ng/ul	0.00
76) Pyrene-d10	19.47	212	954817	31.82	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.35	264	1107028	33.01	ng/ul	0.00

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

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