

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM110717\
 Data File : BM012796.D
 Acq On : 07 Nov 2017 15:20
 Operator : SJ/JU
 Sample : I6164-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ET0D4

Quant Time: Nov 08 00:10:12 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM110417MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 08 00:07:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	89180	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	395150	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.44	164	268615	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	668508	20.00	ng/ul	0.00
77) Chrysene-d12	21.36	240	770318	20.00	ng/ul	0.00
85) Perylene-d12	23.64	264	721923	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	8007	4.85	ng/uL	0.00
5) Phenol-d5	6.97	99	174984	26.37	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	103216	28.70	ng/ul	0.00
9) 2-Chlorophenol-d4	7.34	132	152099	27.52	ng/ul	0.00
13) 4-Methylphenol-d8	8.51	113	152708	26.66	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	78730	26.97	ng/ul	0.00
22) 2-Nitrophenol-d4	9.68	143	93510	27.99	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.22	165	172287	25.16	ng/ul	0.00
29) 4-Chloroaniline-d4	10.73	131	219926	33.28	ng/ul	0.00
43) Dimethylphthalate-d6	13.85	166	635745	28.60	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	689093	25.06	ng/ul	0.00
51) 4-Nitrophenol-d4	14.66	143	54375	15.27	ng/ul	0.00
57) Fluorene-d10	15.43	176	520929	27.36	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.57	200	61919	14.03	ng/ul	0.00
70) Anthracene-d10	17.29	188	849844	26.31	ng/ul	0.00
78) Pyrene-d10	19.57	212	951664	26.62	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.50	264	923546	27.06	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	7.00	94	27636	4.057	ng/ul# 90
44) Dimethylphthalate	13.90	163	315936	14.323	ng/ul 100

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

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