

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_N\DATA\BN032218\
 Data File : BN000551.D
 Acq On : 22 Mar 2018 21:37
 Operator : JU/SJ
 Sample : J1905-17DL2 300X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DAM52DL2

Quant Time: Mar 23 07:58:28 2018
 Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 23 02:22:02 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.41	152	179576	20.00	ng/ul	0.00
18) Naphthalene-d8	11.25	136	764789	20.00	ng/ul	0.00
35) Acenaphthene-d10	15.03	164	430461	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.76	188	930129	20.00	ng/ul	0.00
75) Chrysene-d12	21.88	240	857789	20.00	ng/ul	-0.01
83) Perylene-d12	24.36	264	734730	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	7.56	99	555	0.03	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.71	67	1691	0.17	ng/ul	0.00
9) 2-Chlorophenol-d4	7.94	132	1047	0.08	ng/ul	0.00
13) 4-Methylphenol-d8	9.13	113	932	0.07	ng/ul	0.00
19) Nitrobenzene-d5	9.60	128	579	0.09	ng/ul	0.00
22) 2-Nitrophenol-d4	10.33	143	506	0.08	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.87	165	1121	0.10	ng/ul	0.00
29) 4-Chloroaniline-d4	11.35	131	2392	0.20	ng/ul	-0.04
43) Dimethylphthalate-d6	14.42	166	4281	0.13	ng/ul	0.00
46) Acenaphthylene-d8	14.73	160	5286	0.12	ng/ul	0.00
51) 4-Nitrophenol-d4	15.22	143	1199	0.19	ng/ul	0.00
57) Fluorene-d10	16.01	176	6250	0.22	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.14	200	789	0.15	ng/ul	0.00
70) Anthracene-d10	17.85	188	7953	0.18	ng/ul	0.00
76) Pyrene-d10	20.11	212	5912	0.12	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.20	264	4557	0.13	ng/ul	-0.01

Target Compounds

					Qvalue	
21) Isophorone	10.15	82	42227	1.445	ng/ul#	94
55) 2,3,4,6-Tetrachlorophenol	15.65	232	119540	17.115	ng/ul#	92
68) Pentachlorophenol	17.41	266	993976	188.077	ng/ul	95

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

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