

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032519\
 Data File : BM019422.D
 Acq On : 25 Mar 2019 09:32
 Operator : JU/SJ
 Sample : K2027-08DL 10X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0987DL

Quant Time: Mar 26 06:34:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 25 03:11:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	127956	20.00	ng/ul	0.00
18) Naphthalene-d8	10.39	136	525696	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.26	164	296005	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	640442	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	643834	20.00	ng/ul	0.00
86) Perylene-d12	23.44	264	714079	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	498	0.15	ng/uL	0.03
5) Phenol-d5	6.80	99	6641	0.58	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	6.96	67	21290	2.79	ng/ul	0.00
9) 2-Chlorophenol-d4	7.15	132	22620	2.62	ng/ul	0.00
13) 4-Methylphenol-d8	8.34	113	12051	1.41	ng/ul	0.02
19) Nitrobenzene-d5	8.77	128	12709	3.39	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	13101	3.14	ng/ul	0.01
26) 2,4-Dichlorophenol-d3	10.03	165	21915	2.78	ng/ul	0.02
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	13.69	166	84041	3.52	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	103198	3.46	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.27	176	76382	3.71	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.43	200	3880	0.92	ng/ul	0.02
71) Anthracene-d10	17.12	188	120129	3.91	ng/ul	0.00
79) Pyrene-d10	19.43	212	121900	4.10	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.30	264	135137	3.57	ng/ul	0.00

Target Compounds

				Ovalue	
20) Nitrobenzene	8.82	77	407858	36.537	ng/ul 98
37) 1,2,4,5-Tetrachlorobenzene	12.43	216	44137	4.686	ng/ul# 96
73) 1,2,3,4-Tetrachlorobenzene	13.05	216	161126	17.353	ng/uL 98

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

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