

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032719\
 Data File : BM019533.D
 Acq On : 28 Mar 2019 08:20
 Operator : JU/SJ
 Sample : K2069-10DL 5X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleID :
 C09C9DL

Manual Integrations
 APPROVED

Sohil
 3/28/2019 3:18:42 PM

Quant Time: Apr 09 04:03:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 28 05:07:19 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	107351	20.00	ng/ul	0.01
18) Naphthalene-d8	10.45	136	417138m	20.00	ng/ul	0.06
36) Acenaphthene-d10	14.25	164	224416	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.00	188	481013	20.00	ng/ul	-0.01
78) Chrysene-d12	21.21	240	548341	20.00	ng/ul	-0.01
86) Perylene-d12	23.42	264	769312	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	6.78	99	11207	1.17	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	29554	4.62	ng/ul	0.00
9) 2-Chlorophenol-d4	7.13	132	31474	4.34	ng/ul	0.00
13) 4-Methylphenol-d8	8.32	113	20678	2.88	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	16911	5.68	ng/ul	0.00
22) 2-Nitrophenol-d4	9.48	143	20252	6.12	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.02	165	35862	5.74	ng/ul	0.01
29) 4-Chloroaniline-d4	10.50	131	106	0.01	ng/ul	-0.04
44) Dimethylphthalate-d6	13.67	166	115030	6.35	ng/ul	0.00
47) Acenaphthylene-d8	13.94	160	142284	6.28	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.40	143	671	0.24	ng/ul	-0.10
58) Fluorene-d10	15.25	176	100912	6.47	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.41	200	7582	2.38	ng/ul	0.00
71) Anthracene-d10	17.10	188	144737	6.27	ng/ul	-0.01
79) Pyrene-d10	19.41	212	164491	6.49	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.27	264	245665	6.02	ng/ul	-0.02

Target Compounds

					Ovalue	
28) Naphthalene	10.48	128	120766	5.589	ng/ul	98
34) 2-Methylnaphthalene	12.05	142	46944	3.100	ng/ul	92
37) 1,2,4,5-Tetrachlorobenzene	12.43	216	9759517	1366.674	ng/ul#	96
41) 1,1'-Biphenyl	13.09	154	240958	12.607	ng/ul#	89
50) Acenaphthene	14.31	153	20965	1.326	ng/ul	97
54) Dibenzofuran	14.66	168	53079	2.405	ng/ul	96
59) Fluorene	15.30	166	37063	2.119	ng/ul	96
67) Hexachlorobenzene	16.30	284	96568	15.267	ng/ul#	89
70) Phenanthrene	17.04	178	242912	8.907	ng/ul	99
73) 1,2,3,4-Tetrachlorobenzene	13.06	216	33159674m	4754.892	ng/uL	
74) Pentachlorobenzene	14.57	250	5872501	805.057	ng/uL	98
77) Fluoranthene	19.07	202	97943	3.134	ng/ul#	83
80) Pyrene	19.44	202	65645	1.990	ng/ul#	82
84) Bis(2-ethylhexyl)phthalate	21.12	149	188313	9.947	ng/ul#	95

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

