

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022820\
 Data File : BN009982.D
 Acq On : 28 Feb 2020 10:34
 Operator : CG/JU
 Sample : L1612-19DL 5X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleID :
 DBDM1DL

Manual Integrations
 APPROVED

mohammad
 3/3/2020 4:58:33 PM

Quant Time: Feb 28 14:45:58 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 26 07:56:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	91551	20.00	ng/ul	0.00
18) Naphthalene-d8	10.12	136	396064	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.02	164	253272	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.77	188	550629	20.00	ng/ul	0.00
78) Chrysene-d12	21.00	240	518080	20.00	ng/ul	0.00
86) Perylene-d12	23.10	264	581389	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	6.60	99	7887m	1.01	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	6.75	67	5931	1.10	ng/ul	0.00
9) 2-Chlorophenol-d4	6.93	132	7472	1.27	ng/ul	0.00
13) 4-Methylphenol-d8	8.13	113	4458	0.72	ng/ul	0.03
19) Nitrobenzene-d5	8.55	128	1949	0.67	ng/ul	0.02
22) 2-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	9.86	165	5807m	0.95	ng/ul	0.09
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	13.46	166	27036	1.43	ng/ul	0.01
47) Acenaphthylene-d8	13.71	160	29551	1.33	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
58) Fluorene-d10	15.04	176	26431	1.62	ng/ul	0.01
63) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
71) Anthracene-d10	16.87	188	41739	1.65	ng/ul	0.00
79) Pyrene-d10	19.19	212	49874	1.84	ng/ul	0.00
90) Benzo(a)pyrene-d12	22.97	264	51330	1.77	ng/ul	0.00

Target Compounds

					Ovalue	
70) Phenanthrene	16.82	178	53120	1.690	ng/ul	98
72) Anthracene	16.91	178	1307919	40.714	ng/ul	100
75) Carbazole	17.20	167	147648	5.219	ng/ul#	97
77) Fluoranthene	18.85	202	84722	2.302	ng/ul	98
80) Pyrene	19.22	202	140863	3.771	ng/ul	95
83) Benzo(a)anthracene	20.99	228	112446	3.158	ng/ul	97
85) Chrysene	21.04	228	107493m	3.067	ng/ul	
88) Benzo(b)fluoranthene	22.49	252	539593	14.315	ng/ul	98
89) Benzo(k)fluoranthene	22.53	252	144289	4.043	ng/ul	97
91) Benzo(a)pyrene	23.02	252	247804	7.309	ng/ul	99
92) Indeno(1,2,3-cd)pyrene	25.14	276	189894	4.718	ng/ul	95
93) Dibenzo(a,h)anthracene	25.14	278	44895	1.304	ng/ul#	91
94) Benzo(g,h,i)perylene	25.76	276	115841	3.432	ng/ul	97

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

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