

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN021020\
 Data File : BN009691.D
 Acq On : 10 Feb 2020 18:36
 Operator : CG/JU
 Sample : L1324-14ME
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GAT69ME

Manual Integrations
 APPROVED

mohammad
 2/12/2020 9:51:34 AM

Quant Time: Feb 11 10:54:22 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN020320MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 07 15:21:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.43	152	108133	20.00	ng/ul	0.00
18) Naphthalene-d8	10.17	136	469198	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.06	164	290146	20.00	ng/ul	-0.01
61) Phenanthrene-d10	16.82	188	595683	20.00	ng/ul	0.00
77) Chrysene-d12	21.04	240	568969	20.00	ng/ul	0.00
85) Perylene-d12	23.14	264	573428	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.07	96	6717	2.55	ng/uL	0.00
5) Phenol-d5	6.63	99	174870	18.54	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.79	67	128634	20.31	ng/ul	0.00
9) 2-Chlorophenol-d4	6.97	132	140416	20.23	ng/ul	0.00
13) 4-Methylphenol-d8	8.14	113	147488	19.64	ng/ul	0.00
19) Nitrobenzene-d5	8.58	128	70530	20.21	ng/ul	0.00
22) 2-Nitrophenol-d4	9.29	143	74170	19.05	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.82	165	142039	19.47	ng/ul	0.00
29) 4-Chloroaniline-d4	10.33	131	155745	19.86	ng/ul	0.00
43) Dimethylphthalate-d6	13.49	166	414277	19.56	ng/ul	0.00
46) Acenaphthylene-d8	13.75	160	560066	21.88	ng/ul	0.00
51) 4-Nitrophenol-d4	14.31	143	60127	15.03	ng/ul	0.00
57) Fluorene-d10	15.07	176	405067	21.75	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.23	200	50053	13.45	ng/ul	0.00
70) Anthracene-d10	16.92	188	635653	23.48	ng/ul	0.00
78) Pyrene-d10	19.23	212	663958	22.20	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.02	264	661411	23.43	ng/ul	-0.01

Target Compounds

					Ovalue	
47) Acenaphthylene	13.78	152	710440	25.366	ng/ul	99
58) Fluorene	15.15	166	29649	1.338	ng/ul#	45
69) Phenanthrene	16.86	178	109669	3.228	ng/ul	99
71) Anthracene	16.96	178	372375	10.803	ng/ul	98
76) Fluoranthene	18.89	202	1936378	49.943	ng/ul	99
79) Pyrene	19.26	202	5802559	139.704	ng/ul	96
82) Benzo(a)anthracene	21.02	228	1963092m	49.971	ng/ul	
84) Chrysene	21.07	228	1619647	41.414	ng/ul	97
87) Benzo(b)fluoranthene	22.53	252	1572020	43.817	ng/ul	97
88) Benzo(k)fluoranthene	22.56	252	500670m	14.170	ng/ul	
90) Benzo(a)pyrene	23.06	252	1387635	41.746	ng/ul	96
91) Indeno(1,2,3-cd)pyrene	25.21	276	668359	16.912	ng/ul	94
92) Dibenzo(a,h)anthracene	25.21	278	227793	6.848	ng/ul#	92
93) Benzo(g,h,i)perylene	25.83	276	608091	18.360	ng/ul	94

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

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