

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN020320\
 Data File : BN009564.D
 Acq On : 04 Feb 2020 15:02
 Operator : CG/JU
 Sample : L1302-03
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C5AL0

Quant Time: Feb 04 15:38:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN020320MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 03 16:39:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	126879	20.00	ng/ul	0.00
18) Naphthalene-d8	10.19	136	533510	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.07	164	333329	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.83	188	707812	20.00	ng/ul	0.00
77) Chrysene-d12	21.04	240	689156	20.00	ng/ul	0.00
85) Perylene-d12	23.16	264	741620	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	3805	1.23	ng/uL	0.00
5) Phenol-d5	6.64	99	77444	7.00	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.80	67	58376	7.85	ng/ul	0.00
9) 2-Chlorophenol-d4	6.98	132	63669	7.82	ng/ul	0.00
13) 4-Methylphenol-d8	8.16	113	45633	5.18	ng/ul	0.00
19) Nitrobenzene-d5	8.59	128	30961	7.80	ng/ul	0.00
22) 2-Nitrophenol-d4	9.29	143	33094	7.47	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.83	165	58906	7.10	ng/ul	0.00
29) 4-Chloroaniline-d4	10.35	131	68178	7.65	ng/ul	0.00
43) Dimethylphthalate-d6	13.50	166	207154	8.51	ng/ul	0.00
46) Acenaphthylene-d8	13.76	160	236208	8.03	ng/ul	0.00
51) 4-Nitrophenol-d4	14.32	143	18941	4.12	ng/ul	0.01
57) Fluorene-d10	15.08	176	180493	8.44	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.23	200	17270	3.91	ng/ul	0.00
70) Anthracene-d10	16.93	188	268759	8.36	ng/ul	0.00
78) Pyrene-d10	19.23	212	343826	9.49	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.02	264	352067	9.64	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.66	94	11940	1.010	ng/ul 98
44) Dimethylphthalate	13.55	163	89182	3.454	ng/ul 98
69) Phenanthrene	16.87	178	83612	2.071	ng/ul 98
76) Fluoranthene	18.90	202	141450	3.070	ng/ul 99
79) Pyrene	19.26	202	132349	2.631	ng/ul 96
82) Benzo(a)anthracene	21.03	228	81227	1.707	ng/ul 99
84) Chrysene	21.08	228	87894	1.855	ng/ul 95
87) Benzo(b)fluoranthene	22.53	252	112090	2.416	ng/ul# 96
90) Benzo(a)pyrene	23.06	252	76688	1.784	ng/ul# 94
91) Indeno(1,2,3-cd)pyrene	25.22	276	54901	1.074	ng/ul 95
93) Benzo(g,h,i)perylene	25.84	276	49107	1.146	ng/ul# 88

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

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