

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022920\
 Data File : BP001976.D
 Acq On : 01 Mar 2020 01:17
 Operator : CG/JU
 Sample : L1732-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0EG6

Quant Time: Mar 02 10:53:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 27 14:59:37 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	155429	20.00	ng/ul	0.02
18) Naphthalene-d8	10.43	136	902275	20.00	ng/ul	0.02
36) Acenaphthene-d10	14.27	164	637785	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1385430	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1388402	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1593503	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.17	96	37331	10.33	ng/uL	0.00
5) Phenol-d5	6.82	99	158214	13.34	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.99	67	434960	69.28	ng/ul	0.01
9) 2-Chlorophenol-d4	7.19	132	555265	54.14	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	253742	26.04	ng/ul	0.00
19) Nitrobenzene-d5	8.79	128	299906	43.27	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	352016	45.14	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.05	165	619336	41.23	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	36003	2.19	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	2042685	42.32	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	2471870	42.82	ng/ul	0.00
52) 4-Nitrophenol-d4	14.48	143	58439	6.33	ng/ul	0.00
58) Fluorene-d10	15.27	176	1806793	43.26	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	267382	35.01	ng/ul	0.00
71) Anthracene-d10	17.13	188	2804454	45.04	ng/ul	0.00
79) Pyrene-d10	19.37	212	3268755	44.97	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	3691349	45.28	ng/ul	0.00

Target Compounds

					Ovalue	
20) Nitrobenzene	8.83	77	2638618	170.857	ng/ul	99
23) 2-Nitrophenol	9.54	139	27629	3.257	ng/ul	98
37) 1,2,4,5-Tetrachlorobenzene	12.46	216	1189011	55.924	ng/ul	99
45) Dimethylphthalate	13.74	163	86845	1.743	ng/ul	98
73) 1,2,3,4-Tetrachlorobenzene	13.07	216	3278527	147.299	ng/uL	99
74) Pentachlorobenzene	14.60	250	93922	4.273	ng/uL	98

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

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