

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN011921\
 Data File : BN013421.D
 Acq On : 19 Jan 2021 10:02
 Operator : CG/JU
 Sample : PB133977BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK977

Quant Time: Jan 19 10:38:08 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN011521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 19 10:08:46 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	467520	20.00	ng/ul	0.00
20) Naphthalene-d8	10.35	136	1891412	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.22	164	1063322	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.97	188	2062713	20.00	ng/ul	0.00
79) Chrysene-d12	21.17	240	1597216	20.00	ng/ul	0.00
88) Perylene-d12	23.36	264	1357040	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.19	96	82021	6.99	ng/uL	0.00
4) Pyridine-d5	3.57	84	1062411	34.90	ng/ul	0.00
7) Phenol-d5	6.76	99	1355635	36.05	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.93	67	818031	37.06	ng/ul	0.00
11) 2-Chlorophenol-d4	7.12	132	1191389	36.88	ng/ul	0.00
15) 4-Methylphenol-d8	8.29	113	1082175	35.83	ng/ul	0.00
21) Nitrobenzene-d5	8.73	128	547251	37.36	ng/ul	0.00
24) 2-Nitrophenol-d4	9.45	143	579181	37.91	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.97	165	1046836	35.49	ng/ul	0.00
31) 4-Chloroaniline-d4	10.49	131	1542864	36.22	ng/ul	0.00
46) Dimethylphthalate-d6	13.64	166	3004640	37.55	ng/ul	0.00
49) Acenaphthylene-d8	13.91	160	3639642	36.28	ng/ul	0.00
54) 4-Nitrophenol-d4	14.42	143	483748	31.59	ng/ul	0.00
60) Fluorene-d10	15.22	176	2582154	36.74	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.34	200	362694	29.79	ng/ul	0.00
73) Anthracene-d10	17.07	188	3780678	37.73	ng/ul	0.00
81) Pyrene-d10	19.37	212	3795762	41.50	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.22	264	3004703	41.78	ng/ul	0.00

Target Compounds Ovalue

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

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