

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN012221\
 Data File : BN013453.D
 Acq On : 21 Jan 2021 14:07
 Operator : CG/JU
 Sample : PB134060BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK060

Quant Time: Jan 21 14:44:12 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN012221.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 21 13:24:42 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	488756	20.00	ng/ul	0.00
20) Naphthalene-d8	10.35	136	1880248	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.22	164	978221	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.97	188	1743552	20.00	ng/ul	0.00
79) Chrysene-d12	21.17	240	1451960	20.00	ng/ul	0.00
88) Perylene-d12	23.36	264	1497665	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.19	96	81321	6.79	ng/uL	0.00
4) Pyridine-d5	3.57	84	1085852	34.70	ng/ul	0.00
7) Phenol-d5	6.76	99	1426712	37.12	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.93	67	830492	36.75	ng/ul	0.00
11) 2-Chlorophenol-d4	7.12	132	1271889	37.83	ng/ul	0.00
15) 4-Methylphenol-d8	8.28	113	1128351	36.99	ng/ul	0.00
21) Nitrobenzene-d5	8.73	128	568924	38.56	ng/ul	0.00
24) 2-Nitrophenol-d4	9.44	143	614843	40.59	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.97	165	1067818	36.82	ng/ul	0.00
31) 4-Chloroaniline-d4	10.48	131	1572995	38.43	ng/ul	0.00
46) Dimethylphthalate-d6	13.63	166	2854146	39.32	ng/ul	0.00
49) Acenaphthylene-d8	13.91	160	3789402	41.04	ng/ul	0.00
54) 4-Nitrophenol-d4	14.42	143	454513	33.44	ng/ul	0.00
60) Fluorene-d10	15.22	176	2383431	37.77	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.33	200	314406	30.95	ng/ul	0.00
73) Anthracene-d10	17.07	188	3391657	40.26	ng/ul	0.00
81) Pyrene-d10	19.37	212	3316511	38.42	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.22	264	3386549	43.35	ng/ul	0.00

Target Compounds Ovalue

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

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