

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011221\
 Data File : BN013392.D
 Acq On : 12 Jan 2021 13:07
 Operator : CG/JU
 Sample : PB134003BL
 Misc : PBBL-ME-S-20
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK003

Quant Time: Jan 12 13:37:57 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN010521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 12 09:49:34 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	358567	20.00	ng/ul	0.00
20) Naphthalene-d8	10.36	136	1476302	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.22	164	864421	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.97	188	1753586	20.00	ng/ul	0.00
79) Chrysene-d12	21.16	240	1712588	20.00	ng/ul	0.00
88) Perylene-d12	23.33	264	1819952	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.20	96	52613	5.59	ng/uL	0.00
4) Pyridine-d5	3.58	84	519295	22.11	ng/ul	0.00
7) Phenol-d5	6.78	99	907635	30.14	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.94	67	529085	31.55	ng/ul	0.00
11) 2-Chlorophenol-d4	7.13	132	800005	30.50	ng/ul	0.00
15) 4-Methylphenol-d8	8.29	113	722583	29.71	ng/ul	0.00
21) Nitrobenzene-d5	8.73	128	354002	30.10	ng/ul	0.00
24) 2-Nitrophenol-d4	9.45	143	367026	30.28	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.98	165	701644	29.30	ng/ul	0.00
31) 4-Chloroaniline-d4	10.49	131	1043583	31.85	ng/ul	0.00
46) Dimethylphthalate-d6	13.64	166	2178035	32.58	ng/ul	0.00
49) Acenaphthylene-d8	13.91	160	2755149	33.40	ng/ul	0.00
54) 4-Nitrophenol-d4	14.42	143	332425	27.22	ng/ul	0.00
60) Fluorene-d10	15.22	176	1863891	32.23	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.33	200	228584	26.48	ng/ul	0.00
73) Anthracene-d10	17.06	188	2889917	33.93	ng/ul	0.00
81) Pyrene-d10	19.36	212	3166330	34.33	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.20	264	3365765	35.42	ng/ul	0.00

Target Compounds Ovalue

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

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