

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN011221\
 Data File : BN013388.D
 Acq On : 12 Jan 2021 09:51
 Operator : CG/JU
 Sample : PB134005BL
 Misc : PBBL-S-06
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK005

Quant Time: Jan 12 10:31:05 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	379959	20.00	ng/ul	0.00
20) Naphthalene-d8	10.36	136	1609247	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.22	164	979383	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.97	188	1992997	20.00	ng/ul	0.00
79) Chrysene-d12	21.16	240	1970952	20.00	ng/ul	0.00
88) Perylene-d12	23.34	264	2116386	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.20	96	55890	5.61	ng/uL	0.00
4) Pyridine-d5	3.59	84	558481	22.44	ng/ul	0.00
7) Phenol-d5	6.78	99	983366	30.82	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.95	67	569722	32.06	ng/ul	0.00
11) 2-Chlorophenol-d4	7.13	132	853243	30.70	ng/ul	0.00
15) 4-Methylphenol-d8	8.29	113	789798	30.64	ng/ul	0.00
21) Nitrobenzene-d5	8.74	128	388941	30.34	ng/ul	0.00
24) 2-Nitrophenol-d4	9.45	143	415518	31.44	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.98	165	771319	29.55	ng/ul	0.00
31) 4-Chloroaniline-d4	10.49	131	1177410	32.96	ng/ul	0.00
46) Dimethylphthalate-d6	13.65	166	2482357	32.78	ng/ul	0.00
49) Acenaphthylene-d8	13.92	160	3132709	33.52	ng/ul	0.00
54) 4-Nitrophenol-d4	14.42	143	402530	29.09	ng/ul	0.00
60) Fluorene-d10	15.22	176	2154433	32.88	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.33	200	291862	29.75	ng/ul	0.00
73) Anthracene-d10	17.07	188	3331479	34.42	ng/ul	0.00
81) Pyrene-d10	19.37	212	3667770	34.55	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.20	264	3974529	35.96	ng/ul	0.00

Target Compounds Ovalue

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

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