

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011121\
 Data File : BN013381.D
 Acq On : 11 Jan 2021 13:12
 Operator : CG/JU
 Sample : PB133958BL
 Misc : PBBL-ME-S-04
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK958

Quant Time: Jan 11 13:42:26 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN010521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 08 12:05:50 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	315088	20.00	ng/ul	0.00
20) Naphthalene-d8	10.36	136	1307546	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.22	164	771730	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.97	188	1555095	20.00	ng/ul	0.00
79) Chrysene-d12	21.16	240	1519377	20.00	ng/ul	-0.01
88) Perylene-d12	23.33	264	1611230	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.20	96	46664	5.64	ng/uL	0.00
4) Pyridine-d5	3.59	84	456445	22.12	ng/ul	0.00
7) Phenol-d5	6.78	99	795327	30.06	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.95	67	473194	32.11	ng/ul	0.00
11) 2-Chlorophenol-d4	7.13	132	700462	30.39	ng/ul	0.00
15) 4-Methylphenol-d8	8.30	113	638404	29.87	ng/ul	0.00
21) Nitrobenzene-d5	8.74	128	305124	29.30	ng/ul	0.00
24) 2-Nitrophenol-d4	9.45	143	324266	30.20	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.98	165	613821	28.94	ng/ul	0.00
31) 4-Chloroaniline-d4	10.50	131	933223	32.16	ng/ul	0.00
46) Dimethylphthalate-d6	13.65	166	1906913	31.95	ng/ul	0.00
49) Acenaphthylene-d8	13.92	160	2442336	33.17	ng/ul	0.00
54) 4-Nitrophenol-d4	14.42	143	289491	26.55	ng/ul	0.00
60) Fluorene-d10	15.22	176	1674696	32.43	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.33	200	201258	26.29	ng/ul	0.00
73) Anthracene-d10	17.07	188	2600652	34.43	ng/ul	0.00
81) Pyrene-d10	19.37	212	2845998	34.78	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.20	264	3039966	36.13	ng/ul	-0.02

Target Compounds Ovalue

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

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