

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN042021\
 Data File : BN014328.D
 Acq On : 19 Apr 2021 17:21
 Operator : CG/JU
 Sample : PB135369BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK369

Quant Time: Apr 19 17:54:31 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN041521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 19 15:27:04 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	154949	20.00	ng/ul	0.00
20) Naphthalene-d8	10.46	136	608848	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.33	164	338835	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.07	188	711199	20.00	ng/ul	0.00
79) Chrysene-d12	21.27	240	736267	20.00	ng/ul	0.00
88) Perylene-d12	23.50	264	793572	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	33882	8.32	ng/uL	0.00
4) Pyridine-d5	3.61	84	433717	40.21	ng/ul	-0.01
7) Phenol-d5	6.85	99	509790	39.35	ng/ul	-0.01
9) Bis-(2-Chloroethyl)ether-d	7.02	67	324472	41.83	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.22	132	428638	42.07	ng/ul	0.00
15) 4-Methylphenol-d8	8.38	113	389040	39.62	ng/ul	-0.01
21) Nitrobenzene-d5	8.83	128	193995	44.09	ng/ul	0.00
24) 2-Nitrophenol-d4	9.55	143	183116	42.50	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.09	165	368354	39.69	ng/ul	0.00
31) 4-Chloroaniline-d4	10.60	131	541682	39.34	ng/ul	0.00
46) Dimethylphthalate-d6	13.74	166	1134480	46.22	ng/ul	0.00
49) Acenaphthylene-d8	14.02	160	1432654	48.51	ng/ul	0.00
54) 4-Nitrophenol-d4	14.52	143	175548	36.87	ng/ul	0.00
60) Fluorene-d10	15.32	176	1010075	46.55	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.44	200	124500	32.71	ng/ul	0.00
73) Anthracene-d10	17.17	188	1585555	47.77	ng/ul	0.00
81) Pyrene-d10	19.47	212	1845248	51.63	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.36	264	2079505	49.40	ng/ul	0.00

Target Compounds Ovalue

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

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