

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030921\
 Data File : BN013892.D
 Acq On : 09 Mar 2021 15:08
 Operator : CG/JU
 Sample : PB134931BL
 Misc : SOM-MDL-MES-BL-05
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK31

Quant Time: Mar 09 15:38:21 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN030421MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 09 09:50:00 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	161858	20.00	ng/ul	0.00
18) Naphthalene-d8	10.50	136	680166	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.35	164	402408	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.09	188	816009	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	777776	20.00	ng/ul	0.00
86) Perylene-d12	23.49	264	797959	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.22	96	25956	6.37	ng/uL	0.00
5) Phenol-d5	6.88	99	458740	35.05	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	293570	36.33	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	369013	36.86	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	355505	35.86	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	167435	36.79	ng/ul	0.00
22) 2-Nitrophenol-d4	9.59	143	165379	38.04	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.12	165	331510	34.90	ng/ul	0.00
29) 4-Chloroaniline-d4	10.63	131	510618	43.42	ng/ul	0.00
44) Dimethylphthalate-d6	13.76	166	1110627	41.10	ng/ul	0.00
47) Acenaphthylene-d8	14.05	160	1379114	38.59	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	168310	33.21	ng/ul	0.00
58) Fluorene-d10	15.35	176	948378	39.75	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.46	200	112240	29.59	ng/ul	0.00
71) Anthracene-d10	17.19	188	1494579	41.91	ng/ul	0.00
79) Pyrene-d10	19.48	212	1680517	44.42	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.36	264	1700364	42.80	ng/ul	0.00

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

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