

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN042021\
 Data File : BN014333.D
 Acq On : 19 Apr 2021 20:57
 Operator : CG/JU
 Sample : K4649-13MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ATCLPMSD

Quant Time: Apr 20 01:57:12 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	154697	20.00	ng/ul	0.00
20) Naphthalene-d8	10.47	136	596221	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.33	164	329097	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.07	188	699258	20.00	ng/ul	0.00
79) Chrysene-d12	21.27	240	708187	20.00	ng/ul	0.00
88) Perylene-d12	23.50	264	745518	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
4) Pyridine-d5	3.63	84	789	0.07	ng/ul	0.00
7) Phenol-d5	0.00	99	0	0.00	ng/ul	
9) Bis-(2-Chloroethyl)ether-d	0.00	67	0	0.00	ng/ul	
11) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
15) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
21) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
24) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
28) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
31) 4-Chloroaniline-d4	10.48	131	587	0.04	ng/ul	-0.12
46) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
49) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
54) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
60) Fluorene-d10	0.00	176	0	0.00	ng/ul	
65) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
73) Anthracene-d10	17.17	188	2665	0.08	ng/ul	0.00
81) Pyrene-d10	0.00	212	0	0.00	ng/ul	
92) Benzo(a)pyrene-d12	23.50	264	736983	18.63	ng/ul	0.14

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

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