

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034092.D
 Acq On : 20 Sep 2024 18:06
 Operator : JU/RC
 Sample : P4003-04
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW17811-20240911

Quant Time: Sep 21 01:19:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 18 16:09:28 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.430	152	6191	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	17100	0.400	ng	0.00
13) Acenaphthene-d10	14.072	164	9999	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	20861	0.400	ng	# 0.00
29) Chrysene-d12	21.053	240	13364	0.400	ng	0.00
35) Perylene-d12	23.180	264	10532	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	1912	0.109	ng	0.00
5) Phenol-d6	6.629	99	1435	0.070	ng	0.00
8) Nitrobenzene-d5	8.568	82	3626	0.277	ng	0.00
11) 2-Methylnaphthalene-d10	11.788	152	7589	0.301	ng	0.00
14) 2,4,6-Tribromophenol	15.579	330	1087	0.240	ng	0.00
15) 2-Fluorobiphenyl	12.693	172	12234	0.301	ng	0.00
27) Fluoranthene-d10	18.880	212	19983	0.380	ng	0.00
31) Terphenyl-d14	19.498	244	12833	0.481	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.075	88	6032	0.779	ng	# 74
6) bis(2-Chloroethyl)ether	6.874	93	1934	0.107	ng	98

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

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