

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034087.D
 Acq On : 20 Sep 2024 15:05
 Operator : JU/RC
 Sample : P3957-17
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE104D1-20240911

Quant Time: Sep 20 16:54:51 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	5598	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	14907	0.400	ng	0.00
13) Acenaphthene-d10	14.083	164	8575	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	17660	0.400	ng	# 0.00
29) Chrysene-d12	21.062	240	11408	0.400	ng	0.00
35) Perylene-d12	23.183	264	9752	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	1939	0.122	ng	0.00
5) Phenol-d6	6.636	99	1032	0.056	ng	0.00
8) Nitrobenzene-d5	8.578	82	3545	0.311	ng	0.00
11) 2-Methylnaphthalene-d10	11.791	152	7810	0.355	ng	0.00
14) 2,4,6-Tribromophenol	15.592	330	970	0.250	ng	0.00
15) 2-Fluorobiphenyl	12.693	172	12503	0.358	ng	0.00
27) Fluoranthene-d10	18.880	212	17471	0.392	ng	0.00
31) Terphenyl-d14	19.498	244	10980	0.482	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.068	88	23395	3.343	ng	# 86
24) Pentachlorophenol	16.498	266	565	0.154	ng	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
Data File : BN034087.D
Acq On : 20 Sep 2024 15:05
Operator : JU/RC
Sample : P3957-17
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
RE104D1-20240911

Quant Time: Sep 20 16:54:51 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

