

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034088.D
 Acq On : 20 Sep 2024 15:41
 Operator : JU/RC
 Sample : P3957-18
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DUPE1-20240911

Quant Time: Sep 20 16:55:09 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	7.431	152	5485	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	15180	0.400	ng	0.00
13) Acenaphthene-d10	14.079	164	9163	0.400	ng	0.00
19) Phenanthrene-d10	16.828	188	19611	0.400	ng	#-0.01
29) Chrysene-d12	21.060	240	14571	0.400	ng	# 0.00
35) Perylene-d12	23.184	264	11887	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	3127	0.201	ng	0.00
5) Phenol-d6	6.629	99	1944	0.107	ng	0.00
8) Nitrobenzene-d5	8.579	82	3703	0.319	ng	0.00
11) 2-Methylnaphthalene-d10	11.788	152	8301	0.370	ng	0.00
14) 2,4,6-Tribromophenol	15.586	330	1193	0.287	ng	0.00
15) 2-Fluorobiphenyl	12.690	172	13075	0.351	ng	-0.01
27) Fluoranthene-d10	18.878	212	20761	0.420	ng	0.00
31) Terphenyl-d14	19.501	244	13927	0.479	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.068	88	21924	3.197	ng	# 86
24) Pentachlorophenol	16.493	266	1752	0.429	ng	99

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
Data File : BN034088.D
Acq On : 20 Sep 2024 15:41
Operator : JU/RC
Sample : P3957-18
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DUPE1-20240911

Quant Time: Sep 20 16:55:09 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

