

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091924\
 Data File : BN034048.D
 Acq On : 19 Sep 2024 11:22
 Operator : JU/RC
 Sample : P3845-08DL2 50X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WPDL2

Quant Time: Sep 19 11:49:12 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	7.438	152	7953	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	22663	0.400	ng	# 0.00
13) Acenaphthene-d10	14.083	164	12115	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	27146	0.400	ng	# 0.00
29) Chrysene-d12	21.062	240	21019	0.400	ng	0.00
35) Perylene-d12	23.192	264	22022	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.076	112	65932	2.921	ng	0.00
5) Phenol-d6	6.629	99	80926	3.083	ng	0.00
8) Nitrobenzene-d5	8.579	82	38831	2.240	ng	0.00
11) 2-Methylnaphthalene-d10	11.776	152	5	0.000	ng	-0.02
14) 2,4,6-Tribromophenol	15.580	330	18465	3.364	ng	0.00
15) 2-Fluorobiphenyl	12.693	172	117412	2.383	ng	0.00
27) Fluoranthene-d10	18.890	212	20	0.000	ng	0.00
31) Terphenyl-d14	19.503	244	108123	2.576	ng	0.00
Target Compounds						
20) 4,6-Dinitro-2-methylph...	15.227	198	12269	2.574	ng	# 48
24) Pentachlorophenol	16.498	266	5598	0.989	ng	100

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091924\
Data File : BN034048.D
Acq On : 19 Sep 2024 11:22
Operator : JU/RC
Sample : P3845-08DL2 50X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
PT-ACIDS-WPDL2

Quant Time: Sep 19 11:49:12 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

