

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012422\
 Data File : BP008899.D
 Acq On : 24 Jan 2022 22:00
 Operator : CG/JU
 Sample : N1217-04 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 2256-2261

Quant Time: Jan 25 04:14:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 14 18:05:00 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	434930	20.000	ng	0.00
21) Naphthalene-d8	10.652	136	1719619	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	1054266	20.000	ng	0.00
64) Phenanthrene-d10	17.275	188	1453779	20.000	ng	# 0.02
76) Chrysene-d12	21.363	240	1737451	20.000	ng	0.02
86) Perylene-d12	23.780	264	1811833	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	21641	0.769	ng	0.00
7) Phenol-d6	7.046	99	24487	0.719	ng	0.01
23) Nitrobenzene-d5	9.022	82	18553	0.613	ng	0.00
42) 2,4,6-Tribromophenol	16.004	330	17042	1.111	ng	0.01
45) 2-Fluorobiphenyl	13.116	172	75157	0.940	ng	0.00
79) Terphenyl-d14	19.839	244	90948	0.884	ng	0.02
Target Compounds						
						Qvalue
37) 2-Methylnaphthalene	12.316	142	172534	2.527	ng	# 96
38) 1-Methylnaphthalene	12.534	142	184515	2.944	ng	# 99
58) Fluorene	15.551	166	242456	3.179	ng	# 57
71) Phenanthrene	17.316	178	1366471	16.694	ng	# 79

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012422\
Data File : BP008899.D
Acq On : 24 Jan 2022 22:00
Operator : CG/JU
Sample : N1217-04 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
2256-2261

Quant Time: Jan 25 04:14:15 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 14 18:05:00 2022
Response via : Initial Calibration

