

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080522\
 Data File : BM036129.D
 Acq On : 05 Aug 2022 21:13
 Operator : CG/JU
 Sample : N3961-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RVE-175

Quant Time: Aug 06 01:13:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 03:36:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	257807	20.00	ng	0.00
21) Naphthalene-d8	10.663	136	1060357	20.00	ng	# 0.00
39) Acenaphthene-d10	14.498	164	623163	20.00	ng	0.00
64) Phenanthrene-d10	17.239	188	1218340	20.00	ng	0.00
76) Chrysene-d12	21.415	240	1081153	20.00	ng	0.00
86) Perylene-d12	23.780	264	1173732	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	1673013	105.21	ng	0.02
7) Phenol-d6	7.069	99	2108618	104.21	ng	0.04
23) Nitrobenzene-d5	9.028	82	1263921	66.72	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	822927	112.57	ng	0.00
45) 2-Fluorobiphenyl	13.127	172	2874999	68.12	ng	0.00
79) Terphenyl-d14	19.862	244	3893335	71.03	ng	0.00
Target Compounds						
78) Pyrene	19.656	202	168225	2.51	ng	Qvalue 99
84) Bis(2-ethylhexyl)phtha...	21.333	149	232039	5.32	ng	# 95
88) Benzo(b)fluoranthene	23.056	252	143208	2.08	ng	# 45

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080522\
Data File : BM036129.D
Acq On : 05 Aug 2022 21:13
Operator : CG/JU
Sample : N3961-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RVE-175

Quant Time: Aug 06 01:13:39 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 05 03:36:25 2022
Response via : Initial Calibration

