

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080222\
 Data File : BG054489.D
 Acq On : 2 Aug 2022 18:56
 Operator : CG/JU
 Sample : N3895-19
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-57

Quant Time: Aug 03 04:14:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 28 04:30:50 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.987	152	20039	20.000	ng	0.00
21) Naphthalene-d8	10.796	136	76669	20.000	ng	# 0.00
39) Acenaphthene-d10	14.626	164	57043	20.000	ng	0.00
64) Phenanthrene-d10	17.370	188	137075	20.000	ng	# 0.00
76) Chrysene-d12	21.636	240	153710	20.000	ng	# 0.00
86) Perylene-d12	24.856	264	145505	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.572	112	100519	104.732	ng	0.01
7) Phenol-d6	7.188	99	134630	102.370	ng	0.02
23) Nitrobenzene-d5	9.150	82	91158	64.748	ng	0.00
42) 2,4,6-Tribromophenol	16.119	330	120553	100.655	ng	0.00
45) 2-Fluorobiphenyl	13.246	172	275675	68.175	ng	0.00
79) Terphenyl-d14	19.979	244	541988	69.950	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.427	202	24428	2.625	ng	94
78) Pyrene	19.785	202	25617	2.905	ng	95
88) Benzo(b)fluoranthene	23.833	252	21609	2.511	ng	# 85

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080222\
Data File : BG054489.D
Acq On : 2 Aug 2022 18:56
Operator : CG/JU
Sample : N3895-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RVE-57

Quant Time: Aug 03 04:14:30 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 28 04:30:50 2022
Response via : Initial Calibration

