

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121521\
 Data File : BP008422.D
 Acq On : 15 Dec 2021 18:08
 Operator : CG/JU
 Sample : M5017-09
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP1

Quant Time: Dec 15 18:55:50 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 14 17:59:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.740	152	50636	20.000	ng	0.00
21) Naphthalene-d8	10.540	136	178259	20.000	ng	0.00
39) Acenaphthene-d10	14.398	164	113430	20.000	ng	0.00
64) Phenanthrene-d10	17.163	188	240508	20.000	ng	0.00
76) Chrysene-d12	21.263	240	303506	20.000	ng	0.00
86) Perylene-d12	23.633	264	360457	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.346	112	322875	100.698	ng	0.00
7) Phenol-d6	6.940	99	398294	92.490	ng	0.00
23) Nitrobenzene-d5	8.910	82	262964	62.496	ng	0.00
42) 2,4,6-Tribromophenol	15.904	330	156421	93.837	ng	0.00
45) 2-Fluorobiphenyl	13.016	172	512875	57.477	ng	0.00
79) Terphenyl-d14	19.733	244	870039	49.890	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.204	178	56288	4.320	ng	99
75) Fluoranthene	19.180	202	234361	15.933	ng	98
78) Pyrene	19.533	202	180422	8.309	ng	98
83) Chrysene	21.304	228	84154	4.331	ng	98
88) Benzo(b)fluoranthene	22.916	252	80290	3.703	ng	97

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121521\
Data File : BP008422.D
Acq On : 15 Dec 2021 18:08
Operator : CG/JU
Sample : M5017-09
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP1

Quant Time: Dec 15 18:55:50 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 14 17:59:47 2021
Response via : Initial Calibration

