

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111222\
 Data File : BP012631.D
 Acq On : 12 Nov 2022 21:56
 Operator : CG/JU
 Sample : N5509-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-11

Quant Time: Nov 14 03:22:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Nov 13 22:44:57 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	317554	20.000	ng	0.00
21) Naphthalene-d8	10.510	136	1308216	20.000	ng	0.00
39) Acenaphthene-d10	14.369	164	866550	20.000	ng	-0.01
64) Phenanthrene-d10	17.133	188	1856935	20.000	ng	-0.01
76) Chrysene-d12	21.233	240	1664592	20.000	ng	-0.01
86) Perylene-d12	23.574	264	1486098	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.305	112	2999500	175.884	ng	0.00
7) Phenol-d6	6.905	99	3664231	154.794	ng	0.00
23) Nitrobenzene-d5	8.887	82	2593814	109.998	ng	0.00
42) 2,4,6-Tribromophenol	15.875	330	2277346	192.512	ng	-0.01
45) 2-Fluorobiphenyl	12.987	172	5705441	99.952	ng	-0.01
79) Terphenyl-d14	19.710	244	8792694	107.013	ng	0.00
Target Compounds						
31) Naphthalene	10.557	128	429074	6.037	ng	99
32) Benzoic acid	9.910	122	587750	34.131	ng	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111222\
Data File : BP012631.D
Acq On : 12 Nov 2022 21:56
Operator : CG/JU
Sample : N5509-11
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP-11

Quant Time: Nov 14 03:22:48 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Nov 13 22:44:57 2022
Response via : Initial Calibration

