

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
 Data File : BF131129.D
 Acq On : 10 Nov 2022 18:33
 Operator : CG\JU
 Sample : N5495-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-32A

Quant Time: Nov 11 00:50:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 10 04:15:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.886	152	76956	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	287987	20.000	ng	-0.01
39) Acenaphthene-d10	9.933	164	148352	20.000	ng	0.00
64) Phenanthrene-d10	11.421	188	246888	20.000	ng	0.00
76) Chrysene-d12	14.074	240	112845	20.000	ng	0.00
86) Perylene-d12	15.568	264	129868	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	450533	99.561	ng	0.00
7) Phenol-d6	6.516	99	568962	96.172	ng	0.00
23) Nitrobenzene-d5	7.451	82	370183	57.997	ng	-0.01
42) 2,4,6-Tribromophenol	10.727	330	136560	89.070	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	684380	64.399	ng	0.00
79) Terphenyl-d14	13.015	244	434053	67.441	ng	0.00
Target Compounds						
32) Benzoic acid	7.904	122	3617m	6.797	ng	Qvalue
87) Indeno(1,2,3-cd)pyrene	17.074	276	3719	4.187	ng	98
92) Benzo(g,h,i)perylene	17.545	276	3432	4.707	ng	# 90

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131129.D
Acq On : 10 Nov 2022 18:33
Operator : CG\JU
Sample : N5495-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-32A

Quant Time: Nov 11 00:50:13 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 10 04:15:17 2022
Response via : Initial Calibration

