

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110722\
 Data File : BF131048.D
 Acq On : 07 Nov 2022 17:43
 Operator : CG\JU
 Sample : N5415-05
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-18

Quant Time: Nov 08 02:17:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 01:40:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	88318	20.000	ng	0.00
21) Naphthalene-d8	8.181	136	352962	20.000	ng	0.00
39) Acenaphthene-d10	9.945	164	192417	20.000	ng	0.00
64) Phenanthrene-d10	11.433	188	342565	20.000	ng	0.00
76) Chrysene-d12	14.080	240	178043	20.000	ng	-0.01
86) Perylene-d12	15.580	264	186414	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	424058	77.424	ng	0.00
7) Phenol-d6	6.528	99	613919	91.834	ng	0.00
23) Nitrobenzene-d5	7.463	82	405601	53.266	ng	0.00
42) 2,4,6-Tribromophenol	10.733	330	152372	71.926	ng	0.00
45) 2-Fluorobiphenyl	9.263	172	725579	51.390	ng	0.00
79) Terphenyl-d14	13.022	244	657005	59.163	ng	-0.01
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.086	276	2884	3.493	ng	Qvalue # 83
92) Benzo(g,h,i)perylene	17.545	276	2975	3.603	ng	# 88

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110722\
Data File : BF131048.D
Acq On : 07 Nov 2022 17:43
Operator : CG\JU
Sample : N5415-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
R-18

Quant Time: Nov 08 02:17:07 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 08 01:40:43 2022
Response via : Initial Calibration

