

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062221\
 Data File : BF124427.D
 Acq On : 22 Jun 2021 23:18
 Operator : JU/CG
 Sample : M2789-23
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1ST-WALLS

Quant Time: Jun 23 02:52:06 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 15 20:09:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	37967	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	133424	20.000	ng	# 0.00
39) Acenaphthene-d10	9.822	164	65291	20.000	ng	0.00
64) Phenanthrene-d10	11.316	188	97438	20.000	ng	0.00
76) Chrysene-d12	13.974	240	76702	20.000	ng	# 0.01
86) Perylene-d12	15.439	264	61390	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	160158	63.501	ng	0.00
7) Phenol-d6	6.434	99	191312	56.432	ng	-0.01
23) Nitrobenzene-d5	7.351	82	150114	44.623	ng	0.00
42) 2,4,6-Tribromophenol	10.621	330	48313	52.229	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	276336	53.259	ng	0.00
79) Terphenyl-d14	12.910	244	226839	40.604	ng	0.00
Target Compounds						
						Qvalue
32) Benzoic acid	8.010	122	8156	6.253	ng	88
71) Phenanthrene	11.339	178	15762	2.590	ng	# 90
75) Fluoranthene	12.533	202	14506	2.270	ng	93
84) Bis(2-ethylhexyl)phtha...	13.945	149	29731	8.365	ng	# 95

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062221\
Data File : BF124427.D
Acq On : 22 Jun 2021 23:18
Operator : JU/CG
Sample : M2789-23
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
1ST-WALLS

Quant Time: Jun 23 02:52:06 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 15 20:09:26 2021
Response via : Initial Calibration

