

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062221\
 Data File : BF124426.D
 Acq On : 22 Jun 2021 22:46
 Operator : JU/CG
 Sample : M2789-31
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BASEMENT-WALLS

Quant Time: Jun 23 02:51:57 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	38566	20.000	ng	0.00
21) Naphthalene-d8	8.063	136	140841	20.000	ng	#-0.01
39) Acenaphthene-d10	9.822	164	74772	20.000	ng	0.00
64) Phenanthrene-d10	11.310	188	101708	20.000	ng	0.00
76) Chrysene-d12	13.963	240	81921	20.000	ng	0.00
86) Perylene-d12	15.427	264	72455	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	107802	42.078	ng	0.00
7) Phenol-d6	6.428	99	208066	60.420	ng	-0.02
23) Nitrobenzene-d5	7.351	82	203766	57.381	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	19941	18.824	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	399327	67.204	ng	0.00
79) Terphenyl-d14	12.898	244	297266	49.820	ng	0.00
Target Compounds						
71) Phenanthrene	11.333	178	13014	2.048	ng	98
84) Bis(2-ethylhexyl)phtha...	13.939	149	147078	38.746	ng	97

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

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