

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090622\
 Data File : BF130148.D
 Acq On : 06 Sep 2022 20:52
 Operator : CG\JU
 Sample : N4501-14
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-15-COMP

Quant Time: Sep 07 02:18:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 07 02:05:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.698	152	214105	20.000	ng	0.00
21) Naphthalene-d8	7.975	136	848919	20.000	ng	0.00
39) Acenaphthene-d10	9.722	164	459978	20.000	ng	-0.01
64) Phenanthrene-d10	11.204	188	799999	20.000	ng	0.00
76) Chrysene-d12	13.833	240	496387	20.000	ng	-0.01
86) Perylene-d12	15.233	264	372649	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.310	112	1765904	140.020	ng	0.01
7) Phenol-d6	6.334	99	2303411	146.497	ng	0.00
23) Nitrobenzene-d5	7.263	82	1352396	102.381	ng	0.00
42) 2,4,6-Tribromophenol	10.516	330	685081	163.618	ng	0.00
45) 2-Fluorobiphenyl	9.057	172	2454199	88.460	ng	0.00
79) Terphenyl-d14	12.798	244	2679236	93.599	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	11.227	178	181089	4.174	ng	99
75) Fluoranthene	12.410	202	195773	4.551	ng	98
78) Pyrene	12.633	202	165831	3.725	ng	99

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

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