

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF032122\
 Data File : BF127445.D
 Acq On : 22 Mar 2022 09:56
 Operator : CG\JU
 Sample : N1962-16
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C-19-18-19

Quant Time: Mar 22 15:23:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF032122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 22 02:16:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.846	152	100931	20.000	ng	0.00
21) Naphthalene-d8	8.128	136	373983	20.000	ng	0.00
39) Acenaphthene-d10	9.887	164	222892	20.000	ng	0.00
64) Phenanthrene-d10	11.375	188	409437	20.000	ng	0.00
76) Chrysene-d12	14.022	240	287951	20.000	ng	0.00
86) Perylene-d12	15.504	264	202575	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	756493	119.639	ng	0.02
7) Phenol-d6	6.487	99	944921	109.325	ng	0.00
23) Nitrobenzene-d5	7.416	82	766487	87.608	ng	0.00
42) 2,4,6-Tribromophenol	10.681	330	314706	132.543	ng	0.00
45) 2-Fluorobiphenyl	9.204	172	1300708	85.213	ng	0.00
79) Terphenyl-d14	12.969	244	1567827	89.132	ng	0.00
Target Compounds						
31) Naphthalene	8.151	128	127671	6.430	ng	99
71) Phenanthrene	11.398	178	57110	2.675	ng	99
73) Carbazole	11.616	167	80656	3.982	ng	98

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

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