

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF040622\
 Data File : BF127667.D
 Acq On : 07 Apr 2022 00:28
 Operator : CG\JU
 Sample : N2267-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LINDEN-SOUTH

Quant Time: Apr 07 02:41:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 06 13:55:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.963	152	101419	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	380019	20.000	ng	0.00
39) Acenaphthene-d10	10.010	164	205569	20.000	ng	0.00
64) Phenanthrene-d10	11.498	188	418211	20.000	ng	0.00
76) Chrysene-d12	14.151	240	367215	20.000	ng	0.00
86) Perylene-d12	15.674	264	178482	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.592	112	805082	128.988	ng	0.01
7) Phenol-d6	6.581	99	913780	111.911	ng	0.00
23) Nitrobenzene-d5	7.528	82	788263	92.474	ng	0.00
42) 2,4,6-Tribromophenol	10.798	330	332263	151.887	ng	0.00
45) 2-Fluorobiphenyl	9.327	172	1387187	95.462	ng	0.00
79) Terphenyl-d14	13.098	244	1853427	83.310	ng	0.00

Target Compounds	Qvalue

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

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