

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF040122\
 Data File : BF127610.D
 Acq On : 01 Apr 2022 14:11
 Operator : CG\JU
 Sample : N2217-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-1D

Quant Time: Apr 01 16:13:51 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 29 07:13:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	106092	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	398088	20.000	ng	0.00
39) Acenaphthene-d10	9.869	164	236968	20.000	ng	0.00
64) Phenanthrene-d10	11.363	188	428472	20.000	ng	0.00
76) Chrysene-d12	14.015	240	285477	20.000	ng	# 0.00
86) Perylene-d12	15.498	264	200159	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	783244	117.844	ng	0.01
7) Phenol-d6	6.481	99	1076963	118.541	ng	0.00
23) Nitrobenzene-d5	7.404	82	750335	80.569	ng	0.00
42) 2,4,6-Tribromophenol	10.669	330	297991	118.048	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	1260354	77.665	ng	0.00
79) Terphenyl-d14	12.951	244	1460930	83.775	ng	0.00

Target Compounds	Qvalue

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

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