

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090922\
 Data File : BF130196.D
 Acq On : 09 Sep 2022 15:46
 Operator : CG\JU
 Sample : N4537-02DL 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW2DL

Quant Time: Sep 09 23:47:11 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	6.692	152	187741	20.000	ng	0.00
21) Naphthalene-d8	7.975	136	714652	20.000	ng	0.00
39) Acenaphthene-d10	9.722	164	359037	20.000	ng	-0.01
64) Phenanthrene-d10	11.204	188	529197	20.000	ng	0.00
76) Chrysene-d12	13.833	240	342415	20.000	ng	-0.01
86) Perylene-d12	15.227	264	396896	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.287	112	257503	23.285	ng	-0.01
7) Phenol-d6	6.316	99	185946	13.487	ng	-0.02
23) Nitrobenzene-d5	7.257	82	454279	40.852	ng	-0.01
42) 2,4,6-Tribromophenol	10.510	330	171788	52.563	ng	-0.01
45) 2-Fluorobiphenyl	9.051	172	859881	39.707	ng	0.00
79) Terphenyl-d14	12.786	244	618744	31.336	ng	-0.01
Target Compounds						
20) 3+4-Methylphenols	7.104	107	673846	56.268	ng	Qvalue 86

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

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