

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090722\
 Data File : BF130167.D
 Acq On : 07 Sep 2022 17:53
 Operator : CG\JU
 Sample : N4495-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-345

Quant Time: Sep 08 02:52:22 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	175636	20.000	ng	0.00
21) Naphthalene-d8	7.975	136	687421	20.000	ng	0.00
39) Acenaphthene-d10	9.722	164	349909	20.000	ng	-0.01
64) Phenanthrene-d10	11.204	188	500160	20.000	ng	0.00
76) Chrysene-d12	13.833	240	318462	20.000	ng	-0.01
86) Perylene-d12	15.233	264	360960	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.310	112	1461899	141.303	ng	0.01
7) Phenol-d6	6.328	99	1671966	129.627	ng	0.00
23) Nitrobenzene-d5	7.263	82	1141144	106.684	ng	0.00
42) 2,4,6-Tribromophenol	10.510	330	447967	140.642	ng	-0.01
45) 2-Fluorobiphenyl	9.051	172	2007028	95.098	ng	0.00
79) Terphenyl-d14	12.792	244	1644941	89.572	ng	0.00
Target Compounds						
32) Benzoic acid	7.686	122	23701	9.330	ng	Qvalue 90

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

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