

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111720\
 Data File : BF122359.D
 Acq On : 17 Nov 2020 22:00
 Operator : JU/CG
 Sample : L4769-10
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-359

Quant Time: Nov 18 00:24:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 12:25:42 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.863	152	263110	20.00	ng	0.00
21) Naphthalene-d8	8.151	136	1008597	20.00	ng	0.00
39) Acenaphthene-d10	9.904	164	538702	20.00	ng	0.00
64) Phenanthrene-d10	11.392	188	903922	20.00	ng	0.00
76) Chrysene-d12	14.027	240	691655	20.00	ng	0.00
86) Perylene-d12	15.498	264	707782	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	53748	3.14	ng	0.00
7) Phenol-d6	6.481	99	873220	38.95	ng	-0.01
23) Nitrobenzene-d5	7.428	82	1080407	65.58	ng	0.00
42) 2,4,6-Tribromophenol	10.692	330	4136	0.72	ng	0.00
45) 2-Fluorobiphenyl	9.227	172	1971550	57.18	ng	0.00
79) Terphenyl-d14	12.974	244	1824426	55.88	ng	0.00
Target Compounds						
15) Benzyl Alcohol	6.998	79	33260	2.09	ng	96
50) Dimethylphthalate	9.616	163	325383	8.50	ng	99
80) Butylbenzylphthalate	13.451	149	175504	11.68	ng	96

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

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