

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071222\
 Data File : BF129193.D
 Acq On : 12 Jul 2022 17:13
 Operator : CG\JU
 Sample : N3648-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CMR-TP2-0708

Quant Time: Jul 12 18:39:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 12 14:13:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.933	152	335392	20.000	ng	0.00
21) Naphthalene-d8	8.216	136	1281609	20.000	ng	0.00
39) Acenaphthene-d10	9.974	164	453537	20.000	ng	0.00
64) Phenanthrene-d10	11.468	188	763645	20.000	ng	0.00
76) Chrysene-d12	14.109	240	446996	20.000	ng	-0.01
86) Perylene-d12	15.621	264	420881	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.580	112	2686478	135.419	ng	0.02
7) Phenol-d6	6.563	99	3151965	127.929	ng	0.00
23) Nitrobenzene-d5	7.498	82	2365862	110.108	ng	0.00
42) 2,4,6-Tribromophenol	10.768	330	861979	174.789	ng	0.00
45) 2-Fluorobiphenyl	9.292	172	2968563	104.712	ng	0.00
79) Terphenyl-d14	13.056	244	2942465	136.699	ng	0.00
Target Compounds						
32) Benzoic acid	8.004	122	497638	35.557	ng	Qvalue 94

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

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