

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071222\
 Data File : BF129190.D
 Acq On : 12 Jul 2022 15:37
 Operator : CG\JU
 Sample : N3648-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CMR-TP1-0708

Quant Time: Jul 12 17:16:04 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	6.934	152	324963	20.000	ng	0.00
21) Naphthalene-d8	8.216	136	1237380	20.000	ng	# 0.00
39) Acenaphthene-d10	9.975	164	428232	20.000	ng	0.00
64) Phenanthrene-d10	11.463	188	719939	20.000	ng	# 0.00
76) Chrysene-d12	14.110	240	405312	20.000	ng	-0.01
86) Perylene-d12	15.621	264	384619	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	2353390	122.436	ng	0.02
7) Phenol-d6	6.563	99	2846870	119.254	ng	0.00
23) Nitrobenzene-d5	7.498	82	2278618	109.838	ng	0.00
42) 2,4,6-Tribromophenol	10.769	330	825730	177.333	ng	0.00
45) 2-Fluorobiphenyl	9.292	172	2842495	106.190	ng	0.00
79) Terphenyl-d14	13.057	244	2840290	145.523	ng	0.00
Target Compounds						
32) Benzoic acid	7.969	122	309151	22.879	ng	Qvalue 99

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

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