

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112222\
 Data File : BG055736.D
 Acq On : 22 Nov 2022 12:12
 Operator : CG/JU
 Sample : N5598-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FRAC-TANK-0760533

Quant Time: Nov 23 02:00:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 23 01:35:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.244	152	51617	20.000	ng	0.00
21) Naphthalene-d8	11.076	136	215475	20.000	ng	0.00
39) Acenaphthene-d10	14.866	164	143537	20.000	ng	0.00
64) Phenanthrene-d10	17.610	188	304679	20.000	ng	0.00
76) Chrysene-d12	21.905	240	283301	20.000	ng	0.00
86) Perylene-d12	25.312	264	283011	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.777	112	362566	126.963	ng	0.00
7) Phenol-d6	7.392	99	435282	114.507	ng	0.00
23) Nitrobenzene-d5	9.419	82	333585	81.827	ng	0.00
42) 2,4,6-Tribromophenol	16.352	330	246856	122.134	ng	0.00
45) 2-Fluorobiphenyl	13.491	172	796720	83.155	ng	0.00
79) Terphenyl-d14	20.195	244	1197256	84.527	ng	0.00

Target Compounds	Qvalue

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

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