

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110322\
 Data File : BG055445.D
 Acq On : 3 Nov 2022 20:08
 Operator : CG/JU
 Sample : N5306-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20221173-COMP-08-MODIFIED-ELUTRIATE

Quant Time: Nov 04 05:31:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 04 05:28:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.232	152	27147	20.000	ng	0.00
21) Naphthalene-d8	11.064	136	107761	20.000	ng	0.00
39) Acenaphthene-d10	14.848	164	76881	20.000	ng	0.00
64) Phenanthrene-d10	17.586	188	186390	20.000	ng	0.00
76) Chrysene-d12	21.858	240	199037	20.000	ng	0.00
86) Perylene-d12	25.224	264	216406	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.759	112	112511	74.429	ng	0.00
7) Phenol-d6	7.380	99	110451	52.689	ng	0.00
23) Nitrobenzene-d5	9.407	82	180953	89.493	ng	0.00
42) 2,4,6-Tribromophenol	16.335	330	157002	149.715	ng	0.00
45) 2-Fluorobiphenyl	13.479	172	466023	89.873	ng	0.00
79) Terphenyl-d14	20.160	244	812691	79.398	ng	0.00

Target Compounds	Qvalue

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

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