

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042122\
 Data File : BG053239.D
 Acq On : 21 Apr 2022 13:31
 Operator : CG/JU
 Sample : N2434-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RCA-EDISON-PILE-1

Quant Time: Apr 22 02:56:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 08 15:31:24 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.113	152	30954	20.000	ng	0.00
21) Naphthalene-d8	10.927	136	115210	20.000	ng	0.00
39) Acenaphthene-d10	14.734	164	86767	20.000	ng	-0.01
64) Phenanthrene-d10	17.477	188	213517	20.000	ng	-0.01
76) Chrysene-d12	21.766	240	230117	20.000	ng	0.00
86) Perylene-d12	25.061	264	224390	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.681	112	222901	123.439	ng	0.00
7) Phenol-d6	7.273	99	294060	105.600	ng	0.00
23) Nitrobenzene-d5	9.277	82	248795	87.683	ng	0.00
42) 2,4,6-Tribromophenol	16.226	330	185001	134.017	ng	0.00
45) 2-Fluorobiphenyl	13.359	172	545289	86.658	ng	0.00
79) Terphenyl-d14	20.086	244	1151139	84.486	ng	0.00
Target Compounds						
77) Benzidine	19.704	184	83551	12.726	ng	Qvalue 99

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

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