

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041522\
 Data File : BG053169.D
 Acq On : 15 Apr 2022 18:01
 Operator : CG/JU
 Sample : N2394-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CS-1-(DUMPSTER-RIGHT)

Quant Time: Apr 18 03:16:23 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	8.116	152	29685	20.000	ng	0.00
21) Naphthalene-d8	10.930	136	110127	20.000	ng	0.00
39) Acenaphthene-d10	14.737	164	82480	20.000	ng	-0.01
64) Phenanthrene-d10	17.486	188	211212	20.000	ng	0.00
76) Chrysene-d12	21.763	240	244934	20.000	ng	-0.01
86) Perylene-d12	25.064	264	246967	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.684	112	227089	131.134	ng	0.00
7) Phenol-d6	7.276	99	294306	110.206	ng	0.00
23) Nitrobenzene-d5	9.280	82	252346	93.039	ng	0.00
42) 2,4,6-Tribromophenol	16.229	330	191831	146.188	ng	0.00
45) 2-Fluorobiphenyl	13.362	172	544261	90.990	ng	0.00
79) Terphenyl-d14	20.088	244	1288067	88.817	ng	0.00

Target Compounds	Qvalue

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

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