

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041322\
 Data File : BG053138.D
 Acq On : 13 Apr 2022 14:48
 Operator : CG/JU
 Sample : N2368-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E14ST-EB3-041122-5-10

Quant Time: Apr 14 02:56:48 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	8.119	152	30702	20.000	ng	0.00
21) Naphthalene-d8	10.933	136	127933	20.000	ng	0.00
39) Acenaphthene-d10	14.740	164	97506	20.000	ng	0.00
64) Phenanthrene-d10	17.483	188	231758	20.000	ng	0.00
76) Chrysene-d12	21.766	240	231111	20.000	ng	0.00
86) Perylene-d12	25.061	264	239078	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.681	112	202882	113.275	ng	0.00
7) Phenol-d6	7.279	99	301792	109.266	ng	0.00
23) Nitrobenzene-d5	9.277	82	218019	69.195	ng	0.00
42) 2,4,6-Tribromophenol	16.226	330	169664	109.370	ng	0.00
45) 2-Fluorobiphenyl	13.365	172	481413	68.080	ng	0.00
79) Terphenyl-d14	20.085	244	884277	64.621	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.530	178	32598	2.575	ng	99
75) Fluoranthene	19.527	202	43973	2.710	ng	97
78) Pyrene	19.892	202	41544	2.524	ng	96

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

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