

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081123\
 Data File : BG058707.D
 Acq On : 11 Aug 2023 18:02
 Operator : MA/JU
 Sample : 03958-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NEVINS-SB02-Z106-107

Quant Time: Aug 11 23:01:30 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG072823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 28 22:29:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.812	152	44222	20.000	ng	0.00
21) Naphthalene-d8	10.614	136	192084	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	136429	20.000	ng	0.00
64) Phenanthrene-d10	17.206	188	336625	20.000	ng	0.00
76) Chrysene-d12	21.448	240	297118	20.000	ng	0.00
86) Perylene-d12	24.521	264	337368	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.385	112	267007	92.286	ng	0.00
7) Phenol-d6	6.983	99	377098	93.668	ng	0.00
23) Nitrobenzene-d5	8.975	82	225161	57.585	ng	0.00
42) 2,4,6-Tribromophenol	15.949	330	184595	82.550	ng	0.00
45) 2-Fluorobiphenyl	13.076	172	523801	57.537	ng	0.00
79) Terphenyl-d14	19.821	244	965470	61.107	ng	0.00
Target Compounds						
71) Phenanthrene	17.248	178	56218	3.521	ng	98
78) Pyrene	19.627	202	46132	2.281	ng	98

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

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