

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072922\
 Data File : BG054449.D
 Acq On : 29 Jul 2022 14:47
 Operator : CG/JU
 Sample : N3893-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-3

Manual Integrations
 APPROVED

Reviewed By : Christian
 Giraldo 08/01/2022
 Supervised By : Jaqrut
 Upadhyay 08/01/2022

Quant Time: Jul 30 03:18:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 28 04:30:50 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.987	152	17079	20.000	ng	0.00
21) Naphthalene-d8	10.795	136	67634	20.000	ng	# 0.00
39) Acenaphthene-d10	14.620	164	53279	20.000	ng	0.00
64) Phenanthrene-d10	17.370	188	127323	20.000	ng	# 0.00
76) Chrysene-d12	21.635	240	142947	20.000	ng	# 0.00
86) Perylene-d12	24.838	264	152400	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.566	112	100135	122.414	ng	0.00
7) Phenol-d6	7.176	99	135188	120.610	ng	0.00
23) Nitrobenzene-d5	9.150	82	88902	71.581	ng	0.00
42) 2,4,6-Tribromophenol	16.118	330	122837	109.807	ng	0.00
45) 2-Fluorobiphenyl	13.245	172	276959	73.332	ng	0.00
79) Terphenyl-d14	19.979	244	556851	77.279	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.420	202	31251	3.616	ng	95
78) Pyrene	19.785	202	32580	3.972	ng	95
88) Benzo(b)fluoranthene	23.821	252	31937	3.544	ng	98
90) Benzo(a)pyrene	24.685	252	24492	3.182	ng	96

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

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