

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080222\
 Data File : BG054485.D
 Acq On : 2 Aug 2022 16:08
 Operator : CG/JU
 Sample : N3895-12
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-50

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 08/03/2022
 Supervised By :Jagrut Upadhyay 08/04/2022

Quant Time: Aug 03 04:13:59 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	7.984	152	18412	20.000	ng	0.00
21) Naphthalene-d8	10.793	136	70097	20.000	ng	# 0.00
39) Acenaphthene-d10	14.618	164	52423	20.000	ng	0.00
64) Phenanthrene-d10	17.368	188	128302	20.000	ng	# 0.00
76) Chrysene-d12	21.633	240	144552	20.000	ng	# 0.00
86) Perylene-d12	24.847	264	154407	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.570	112	119981	136.056	ng	0.00
7) Phenol-d6	7.185	99	162722	134.664	ng	0.01
23) Nitrobenzene-d5	9.148	82	110834	86.105	ng	0.00
42) 2,4,6-Tribromophenol	16.116	330	153652	139.596	ng	0.00
45) 2-Fluorobiphenyl	13.243	172	341629	91.932	ng	0.00
79) Terphenyl-d14	19.976	244	687437	94.342	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	19.424	202	31738	3.644	ng	98
78) Pyrene	19.782	202	37276	4.494	ng	98
83) Chrysene	21.680	228	23468m	2.707	ng	
88) Benzo(b)fluoranthene	23.831	252	30195	3.307	ng	# 96
90) Benzo(a)pyrene	24.694	252	20704	2.655	ng	94

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

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