

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080422\
 Data File : BG054523.D
 Acq On : 4 Aug 2022 17:44
 Operator : CG/JU
 Sample : N3930-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-99D

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/05/2022
 Supervised By : Jagrut Upadhyay 08/05/2022

Quant Time: Aug 04 18:32:11 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	21248	20.00	ng	0.00
21) Naphthalene-d8	10.792	136	80704	20.00	ng	# 0.00
39) Acenaphthene-d10	14.623	164	57750	20.00	ng	0.00
64) Phenanthrene-d10	17.373	188	131425	20.00	ng	# 0.00
76) Chrysene-d12	21.638	240	142670	20.00	ng	# 0.00
86) Perylene-d12	24.852	264	138033	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.569	112	83964	82.51	ng	0.00
7) Phenol-d6	7.185	99	111931	80.27	ng	0.01
23) Nitrobenzene-d5	9.147	82	77215	52.10	ng	0.00
42) 2,4,6-Tribromophenol	16.121	330	86527	71.36	ng	0.00
45) 2-Fluorobiphenyl	13.242	172	213449	52.14	ng	0.00
79) Terphenyl-d14	19.976	244	357582	49.72	ng	0.00
Target Compounds						
78) Pyrene	19.788	202	25498m	3.11	ng	Qvalue
88) Benzo(b)fluoranthene	23.842	252	18486	2.26	ng	# 72

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

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