

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080522\
 Data File : BM036134.D
 Acq On : 06 Aug 2022 00:17
 Operator : CG/JU
 Sample : N3961-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RVE-179

Manual Integrations
 APPROVED

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

Quant Time: Aug 06 01:14:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 03:36:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	278682	20.00	ng	0.00
21) Naphthalene-d8	10.663	136	1125051	20.00	ng	# 0.00
39) Acenaphthene-d10	14.498	164	641867	20.00	ng	0.00
64) Phenanthrene-d10	17.239	188	1260467	20.00	ng	0.00
76) Chrysene-d12	21.415	240	1176466	20.00	ng	0.00
86) Perylene-d12	23.774	264	1227027	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	1486351	86.47	ng	0.02
7) Phenol-d6	7.069	99	1833195	83.81	ng	0.04
23) Nitrobenzene-d5	9.028	82	1108223	55.14	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	685323	91.01	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	2506398	57.65	ng	0.00
79) Terphenyl-d14	19.862	244	3551463	59.54	ng	0.00
Target Compounds						
78) Pyrene	19.656	202	224084	3.07	ng	Qvalue 99
88) Benzo(b)fluoranthene	23.056	252	162897m	2.27	ng	
90) Benzo(a)pyrene	23.668	252	130373	2.16	ng	# 90

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

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