

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080222\
 Data File : BM036063.D
 Acq On : 02 Aug 2022 22:06
 Operator : CG/JU
 Sample : N3927-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RVE-82

Manual Integrations
 APPROVED

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

Quant Time: Aug 03 00:54:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 17:50:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	363453	20.000	ng	0.00
21) Naphthalene-d8	10.675	136	1556447	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	938608	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1948221	20.000	ng	0.00
76) Chrysene-d12	21.421	240	1801480	20.000	ng	0.00
86) Perylene-d12	23.786	264	1429628	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	2421482	108.015	ng	0.00
7) Phenol-d6	7.040	99	3132928	109.830	ng	0.00
23) Nitrobenzene-d5	9.040	82	1903248	68.449	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	1178164	106.998	ng	0.00
45) 2-Fluorobiphenyl	13.139	172	4297077	67.595	ng	0.00
79) Terphenyl-d14	19.874	244	6666875	72.996	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	17.286	178	337943	3.405	ng	99
75) Fluoranthene	19.304	202	419194	3.773	ng	98
78) Pyrene	19.662	202	359690	3.220	ng	100
88) Benzo(b)fluoranthene	23.062	252	204678	2.444	ng	# 91

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

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