

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080322\
 Data File : BG054499.D
 Acq On : 3 Aug 2022 14:37
 Operator : CG/JU
 Sample : N3935-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-137

Quant Time: Aug 04 00:43:38 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	7.979	152	17558	20.000	ng	0.00
21) Naphthalene-d8	10.787	136	68847	20.000	ng	# 0.00
39) Acenaphthene-d10	14.618	164	55505	20.000	ng	0.00
64) Phenanthrene-d10	17.368	188	134126	20.000	ng	# 0.00
76) Chrysene-d12	21.633	240	151171	20.000	ng	# 0.00
86) Perylene-d12	24.841	264	158649	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.564	112	114431	136.074	ng	0.00
7) Phenol-d6	7.180	99	156207	135.560	ng	0.00
23) Nitrobenzene-d5	9.148	82	103752	82.066	ng	0.00
42) 2,4,6-Tribromophenol	16.116	330	143823	123.411	ng	0.00
45) 2-Fluorobiphenyl	13.237	172	332922	84.614	ng	0.00
79) Terphenyl-d14	19.977	244	665390	87.318	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.418	202	34624	3.803	ng	94
78) Pyrene	19.783	202	41796	4.818	ng	96
81) Benzo(a)anthracene	21.616	228	19256	2.007	ng	91
83) Chrysene	21.680	228	29989	3.308	ng	97
90) Benzo(a)pyrene	24.689	252	25311	3.159	ng	96

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

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