

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080322\
 Data File : BG054506.D
 Acq On : 3 Aug 2022 19:31
 Operator : CG/JU
 Sample : N3935-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-146

Quant Time: Aug 04 00:44:31 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	16572	20.000	ng	0.00
21) Naphthalene-d8	10.788	136	68451	20.000	ng	# 0.00
39) Acenaphthene-d10	14.618	164	55666	20.000	ng	0.00
64) Phenanthrene-d10	17.362	188	137908	20.000	ng	# 0.00
76) Chrysene-d12	21.628	240	149932	20.000	ng	# 0.00
86) Perylene-d12	24.836	264	156811	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.564	112	119373	150.396	ng	0.00
7) Phenol-d6	7.180	99	167936	154.410	ng	0.00
23) Nitrobenzene-d5	9.148	82	109907	87.438	ng	0.00
42) 2,4,6-Tribromophenol	16.117	330	156249	133.686	ng	0.00
45) 2-Fluorobiphenyl	13.238	172	351129	88.984	ng	0.00
79) Terphenyl-d14	19.971	244	676608	89.524	ng	0.00
Target Compounds						
						Qvalue
68) Hexachlorobenzene	16.681	284	5562	2.511	ng	# 88
75) Fluoranthene	19.419	202	32467	3.468	ng	96
78) Pyrene	19.783	202	37655	4.377	ng	98
83) Chrysene	21.675	228	26777	2.978	ng	# 96
88) Benzo(b)fluoranthene	23.825	252	40057	4.319	ng	97
90) Benzo(a)pyrene	24.683	252	28586	3.609	ng	99

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

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