

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080322\
 Data File : BG054508.D
 Acq On : 3 Aug 2022 20:56
 Operator : CG/JU
 Sample : N3927-10
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-83

Quant Time: Aug 04 00:44:46 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	7.977	152	16869	20.000	ng	-0.01
21) Naphthalene-d8	10.791	136	71261	20.000	ng	# 0.00
39) Acenaphthene-d10	14.616	164	56562	20.000	ng	0.00
64) Phenanthrene-d10	17.366	188	136331	20.000	ng	# 0.00
76) Chrysene-d12	21.631	240	148410	20.000	ng	# 0.00
86) Perylene-d12	24.839	264	155429	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.568	112	95326	117.985	ng	0.00
7) Phenol-d6	7.184	99	131777	119.030	ng	0.01
23) Nitrobenzene-d5	9.146	82	89991	68.770	ng	0.00
42) 2,4,6-Tribromophenol	16.114	330	124321	104.683	ng	0.00
45) 2-Fluorobiphenyl	13.241	172	284476	70.950	ng	0.00
79) Terphenyl-d14	19.974	244	544284	72.755	ng	0.00
Target Compounds						
75) Fluoranthene	19.422	202	26818	2.898	ng	Qvalue 91
78) Pyrene	19.781	202	26653	3.130	ng	98
88) Benzo(b)fluoranthene	23.823	252	21658	2.356	ng	# 87

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

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