

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080922\
 Data File : BG054575.D
 Acq On : 9 Aug 2022 18:38
 Operator : CG/JU
 Sample : N3962-09
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-195

Quant Time: Aug 10 07:13:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 08 14:48:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.970	152	19302	20.000	ng	0.00
21) Naphthalene-d8	10.784	136	73700	20.000	ng	# 0.00
39) Acenaphthene-d10	14.615	164	57533	20.000	ng	0.00
64) Phenanthrene-d10	17.365	188	140105	20.000	ng	# 0.00
76) Chrysene-d12	21.630	240	156262	20.000	ng	# 0.00
86) Perylene-d12	24.844	264	159688	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.561	112	92140	99.667	ng	0.00
7) Phenol-d6	7.177	99	125635	99.178	ng	0.00
23) Nitrobenzene-d5	9.139	82	85897	63.469	ng	0.00
42) 2,4,6-Tribromophenol	16.113	330	112170	92.858	ng	0.00
45) 2-Fluorobiphenyl	13.234	172	262063	64.257	ng	0.00
79) Terphenyl-d14	19.974	244	540818	68.659	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	19.415	202	24872	2.615	ng	96
78) Pyrene	19.780	202	30903	3.447	ng	97
85) Di-n-octyl phthalate	22.700	149	25335	3.756	ng	# 98
88) Benzo(b)fluoranthene	23.822	252	28008	2.966	ng	97
90) Benzo(a)pyrene	24.692	252	18132	2.248	ng	# 94

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

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