

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072922\
 Data File : BG054456.D
 Acq On : 29 Jul 2022 19:39
 Operator : CG/JU
 Sample : N3895-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-39D

Quant Time: Jul 30 03:18:56 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	7.991	152	17974	20.000	ng	# 0.00
21) Naphthalene-d8	10.799	136	68255	20.000	ng	# 0.00
39) Acenaphthene-d10	14.624	164	50514	20.000	ng	0.00
64) Phenanthrene-d10	17.368	188	119564	20.000	ng	# 0.00
76) Chrysene-d12	21.633	240	135453	20.000	ng	# 0.00
86) Perylene-d12	24.847	264	135831	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.570	112	76722	89.121	ng	0.00
7) Phenol-d6	7.180	99	102995	87.313	ng	0.00
23) Nitrobenzene-d5	9.148	82	68714	54.823	ng	0.00
42) 2,4,6-Tribromophenol	16.116	330	89658	84.535	ng	0.00
45) 2-Fluorobiphenyl	13.243	172	209419	58.484	ng	0.00
79) Terphenyl-d14	19.977	244	405328	59.363	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.409	178	26974	4.530	ng	98
75) Fluoranthene	19.424	202	46890	5.777	ng	97
78) Pyrene	19.789	202	82592	10.627	ng	97
81) Benzo(a)anthracene	21.616	228	40030	4.655	ng	# 89
83) Chrysene	21.680	228	48914	6.021	ng	# 92
88) Benzo(b)fluoranthene	23.831	252	39988	4.978	ng	98
90) Benzo(a)pyrene	24.695	252	31823	4.639	ng	# 95

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

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