

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060322\
 Data File : BG053815.D
 Acq On : 3 Jun 2022 16:18
 Operator : CG/JU
 Sample : N3170-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20220424-AMENDED-COMPOSITE-D

Quant Time: Jun 04 03:05:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 02 23:22:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.091	152	30654	20.000	ng	0.00
21) Naphthalene-d8	10.900	136	123870	20.000	ng	0.00
39) Acenaphthene-d10	14.713	164	82367	20.000	ng	0.00
64) Phenanthrene-d10	17.457	188	175282	20.000	ng	0.00
76) Chrysene-d12	21.734	240	178614	20.000	ng	0.00
86) Perylene-d12	25.001	264	183936	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.659	112	19889	11.915	ng	0.00
7) Phenol-d6	7.239	99	95872	42.398	ng	0.00
23) Nitrobenzene-d5	9.249	82	186261	98.744	ng	0.00
42) 2,4,6-Tribromophenol	16.193	330	2426	2.426	ng	0.00
45) 2-Fluorobiphenyl	13.344	172	410962	74.642	ng	0.00
79) Terphenyl-d14	20.065	244	619479	68.998	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.507	202	27797	2.323	ng	96
78) Pyrene	19.866	202	31722	2.793	ng	99
83) Chrysene	21.781	228	23279	2.080	ng	# 94
84) Bis(2-ethylhexyl)phtha...	21.628	149	146083	27.422	ng	97
88) Benzo(b)fluoranthene	23.967	252	24849	2.176	ng	# 97
90) Benzo(a)pyrene	24.842	252	22729	2.353	ng	# 90

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

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