

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042523\
 Data File : BG057185.D
 Acq On : 25 Apr 2023 20:43
 Operator : CG/JU
 Sample : 02470-14
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-11-COMP

Quant Time: Apr 26 00:25:22 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 25 04:59:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.385	152	17690	20.000	ng	0.00
21) Naphthalene-d8	11.222	136	73571	20.000	ng	0.00
39) Acenaphthene-d10	15.000	164	53326	20.000	ng	0.00
64) Phenanthrene-d10	17.737	188	117478	20.000	ng	# 0.00
76) Chrysene-d12	22.049	240	92069	20.000	ng	#-0.01
86) Perylene-d12	25.579	264	100498	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.900	112	69165	70.853	ng	0.00
7) Phenol-d6	7.522	99	108211	76.615	ng	0.00
23) Nitrobenzene-d5	9.560	82	66255	44.520	ng	0.00
42) 2,4,6-Tribromophenol	16.480	330	43534	61.874	ng	0.00
45) 2-Fluorobiphenyl	13.619	172	174216	43.906	ng	-0.01
79) Terphenyl-d14	20.304	244	249791	47.455	ng	-0.01
Target Compounds						
						Qvalue
75) Fluoranthene	19.770	202	22546	3.026	ng	94
78) Pyrene	20.128	202	20907	3.345	ng	97
81) Benzo(a)anthracene	22.025	228	14148	2.235	ng	# 92
83) Chrysene	22.102	228	13185	2.160	ng	# 89
84) Bis(2-ethylhexyl)phtha...	21.879	149	83267	33.438	ng	98
88) Benzo(b)fluoranthene	24.446	252	19631	3.116	ng	# 79
90) Benzo(a)pyrene	25.409	252	12919	2.082	ng	# 81

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

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