

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090822\
 Data File : BG054889.D
 Acq On : 8 Sep 2022 18:36
 Operator : CG/JU
 Sample : N4530-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-18-COMP

Quant Time: Sep 09 02:51:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 09 02:45:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.149	152	40316	20.000	ng	0.00
21) Naphthalene-d8	10.975	136	165152	20.000	ng	0.00
39) Acenaphthene-d10	14.782	164	116192	20.000	ng	0.00
64) Phenanthrene-d10	17.532	188	270585	20.000	ng	0.00
76) Chrysene-d12	21.827	240	256877	20.000	ng	0.00
86) Perylene-d12	25.205	264	268122	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.687	112	226064	97.643	ng	0.00
7) Phenol-d6	7.303	99	312439	98.679	ng	0.00
23) Nitrobenzene-d5	9.324	82	185653	59.013	ng	0.00
42) 2,4,6-Tribromophenol	16.269	330	150510	103.665	ng	0.00
45) 2-Fluorobiphenyl	13.407	172	476483	61.636	ng	0.00
79) Terphenyl-d14	20.129	244	807424	62.280	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.577	202	93633	5.340	ng	98
78) Pyrene	19.941	202	93448	5.217	ng	98
81) Benzo(a)anthracene	21.809	228	61982	3.559	ng	100
83) Chrysene	21.874	228	56136	3.371	ng	99
87) Indeno(1,2,3-cd)pyrene	29.089	276	42166	2.072	ng	95
88) Benzo(b)fluoranthene	24.130	252	76190	4.530	ng	99
90) Benzo(a)pyrene	25.047	252	55460	3.859	ng	98
92) Benzo(g,h,i)perylene	30.317	276	44495	2.656	ng	# 92

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

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