

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG103125\
 Data File : BG064670.D
 Acq On : 1 Nov 2025 00:27
 Operator : RC/JU
 Sample : Q3486-07DL 20X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP6DL

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 11/03/2025
 Supervised By :mohammad ahmed 11/05/2025

Quant Time: Nov 01 01:02:42 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 24 16:40:26 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.219	152	43395	20.000	ng	0.00
21) Naphthalene-d8	11.045	136	171761	20.000	ng	0.00
39) Acenaphthene-d10	14.841	164	128277	20.000	ng	0.00
64) Phenanthrene-d10	17.579	188	266257	20.000	ng	0.00
76) Chrysene-d12	21.856	240	285356	20.000	ng	# 0.00
86) Perylene-d12	25.199	264	320271	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.752	112	9375	3.626	ng	0.00
7) Phenol-d6	7.350	99	13458	3.794	ng	0.00
23) Nitrobenzene-d5	9.377	82	9042	3.162	ng	-0.01
42) 2,4,6-Tribromophenol	16.316	330	6889	3.721	ng	0.00
45) 2-Fluorobiphenyl	13.472	172	21913	2.589	ng	0.00
79) Terphenyl-d14	20.164	244	36277	2.399	ng	-0.01
Target Compounds						Qvalue
49) Acenaphthylene	14.571	152	56044m	5.162	ng	
71) Phenanthrene	17.620	178	199665	13.713	ng	99
75) Fluoranthene	19.618	202	279613	15.329	ng	98
78) Pyrene	19.976	202	590366	31.666	ng	99
81) Benzo(a)anthracene	21.833	228	334740m	17.692	ng	
83) Chrysene	21.903	228	347902	19.324	ng	95
87) Indeno(1,2,3-cd)pyrene	29.036	276	142913	6.066	ng	# 87
88) Benzo(b)fluoranthene	24.136	252	270543	13.776	ng	98
89) Benzo(k)fluoranthene	24.201	252	95925m	4.881	ng	
90) Benzo(a)pyrene	25.047	252	306845	16.997	ng	99
91) Dibenzo(a,h)anthracene	29.083	278	48505m	2.534	ng	
92) Benzo(g,h,i)perylene	29.036	276	142913	6.066	ng	98

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

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