

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092225\
 Data File : BG064426.D
 Acq On : 22 Sep 2025 22:24
 Operator : RC/JU
 Sample : Q3133-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 361

Quant Time: Sep 23 01:54:32 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 28 17:41:58 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.836	152	19413	20.000	ng	-0.02
21) Naphthalene-d8	10.626	136	70387	20.000	ng	#-0.02
39) Acenaphthene-d10	14.463	164	51672	20.000	ng	-0.02
64) Phenanthrene-d10	17.207	188	131581	20.000	ng	#-0.02
76) Chrysene-d12	21.431	240	157394	20.000	ng	#-0.02
86) Perylene-d12	24.428	264	173037	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.433	112	145685	134.638	ng	-0.01
7) Phenol-d6	7.013	99	173387	116.637	ng	-0.02
23) Nitrobenzene-d5	8.993	82	162060	128.406	ng	-0.02
42) 2,4,6-Tribromophenol	15.950	330	152574	222.925	ng	-0.02
45) 2-Fluorobiphenyl	13.088	172	370692	109.500	ng	-0.02
79) Terphenyl-d14	19.827	244	893034	118.377	ng	-0.02

Target Compounds	Qvalue

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

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