

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080525\
 Data File : BF143318.D
 Acq On : 05 Aug 2025 16:24
 Operator : RC/JU
 Sample : Q2759-01DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SOLID-DRUMSDL

Quant Time: Aug 05 17:01:31 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF071725.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 17 15:14:05 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.963	152	98241	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	333512	20.000	ng	0.00
39) Acenaphthene-d10	10.004	164	135009	20.000	ng	0.00
64) Phenanthrene-d10	11.492	188	218932	20.000	ng	0.00
76) Chrysene-d12	14.133	240	198223	20.000	ng	0.01
86) Perylene-d12	15.645	264	138342	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.592	112	128216	20.653	ng	0.00
7) Phenol-d6	6.586	99	160135	20.493	ng	0.00
23) Nitrobenzene-d5	7.516	82	91060	13.610	ng	0.00
42) 2,4,6-Tribromophenol	10.792	330	25650	18.391	ng	0.00
45) 2-Fluorobiphenyl	9.316	172	167510	16.969	ng	0.00
79) Terphenyl-d14	13.068	244	164803	12.372	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.604	94	65165	7.655	ng	96
20) 3+4-Methylphenols	7.363	107	393299	59.246	ng	# 72

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

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