

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062525\
 Data File : BF142845.D
 Acq On : 25 Jun 2025 19:06
 Operator : RC/JU
 Sample : Q2386-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-1-062025

Quant Time: Jun 26 01:19:19 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 24 05:28:21 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	85467	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	320972	20.000	ng	-0.01
39) Acenaphthene-d10	9.916	164	164270	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	221790	20.000	ng	0.00
76) Chrysene-d12	14.051	240	169503	20.000	ng	0.00
86) Perylene-d12	15.539	264	183109	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	542121	105.607	ng	0.02
7) Phenol-d6	6.510	99	577503	95.354	ng	0.00
23) Nitrobenzene-d5	7.439	82	444219	75.912	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	177008	104.711	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	915288	73.196	ng	0.00
79) Terphenyl-d14	12.992	244	564301	45.999	ng	0.00
Target Compounds						
						Qvalue
50) Dimethylphthalate	9.628	163	43526	4.093	ng	98
52) Acenaphthene	9.951	154	75394	7.708	ng	98
55) Dibenzofuran	10.122	168	41605	2.966	ng	98
58) Fluorene	10.463	166	57097	5.213	ng	# 97
71) Phenanthrene	11.427	178	99319	8.267	ng	99
73) Carbazole	11.639	167	29125	2.739	ng	98

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

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