

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061125\
 Data File : BF142738.D
 Acq On : 11 Jun 2025 16:49
 Operator : RC/JU
 Sample : Q2268-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-6-20250605

Quant Time: Jun 11 17:28:39 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061125.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 11 05:56:09 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.893	152	51607	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	177791	20.000	ng	0.00
39) Acenaphthene-d10	9.928	164	91569	20.000	ng	-0.01
64) Phenanthrene-d10	11.416	188	141283	20.000	ng	0.00
76) Chrysene-d12	14.063	240	134836	20.000	ng	0.00
86) Perylene-d12	15.557	264	163138	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	187744	61.960	ng	0.00
7) Phenol-d6	6.516	99	156124	43.781	ng	0.00
23) Nitrobenzene-d5	7.451	82	276551	85.128	ng	0.00
42) 2,4,6-Tribromophenol	10.722	330	121507	121.073	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	576472	83.639	ng	0.00
79) Terphenyl-d14	13.004	244	551740	56.200	ng	0.00
Target Compounds						
10) Phenol	6.528	94	20831	5.255	ng	Qvalue 89
27) 2,4-Dimethylphenol	7.840	122	297906	108.739	ng	# 93

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

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