

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060925\
 Data File : BP024874.D
 Acq On : 09 Jun 2025 12:50
 Operator : RC/JU
 Sample : Q2210-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TW1

Quant Time: Jun 09 13:13:17 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 06 16:20:27 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.607	152	152557	20.000	ng	0.00
21) Naphthalene-d8	10.372	136	617882	20.000	ng	0.00
39) Acenaphthene-d10	14.242	164	395301	20.000	ng	0.00
64) Phenanthrene-d10	17.048	188	788461	20.000	ng	-0.01
76) Chrysene-d12	21.483	240	1056540	20.000	ng	0.00
86) Perylene-d12	24.712	264	1470569	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.237	112	742475	81.238	ng	0.00
7) Phenol-d6	6.813	99	614677	50.831	ng	0.00
23) Nitrobenzene-d5	8.754	82	1169640	91.986	ng	0.00
42) 2,4,6-Tribromophenol	15.766	330	895820	163.912	ng	-0.02
45) 2-Fluorobiphenyl	12.848	172	2487280	84.762	ng	-0.01
79) Terphenyl-d14	19.777	244	4732640	80.282	ng	-0.02

Target Compounds Qvalue

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

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