

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061125\
 Data File : BP024911.D
 Acq On : 11 Jun 2025 14:21
 Operator : RC/JU
 Sample : Q2227-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP07-MHH-WC

Quant Time: Jun 11 14:49:04 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	7.608	152	262982	20.000	ng	0.00
21) Naphthalene-d8	10.384	136	1037736	20.000	ng	0.00
39) Acenaphthene-d10	14.254	164	675868	20.000	ng	0.00
64) Phenanthrene-d10	17.048	188	1424970	20.000	ng	-0.01
76) Chrysene-d12	21.483	240	1742549	20.000	ng	0.00
86) Perylene-d12	24.742	264	2024689	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.237	112	1155321	73.331	ng	0.00
7) Phenol-d6	6.813	99	1520111	72.923	ng	0.00
23) Nitrobenzene-d5	8.760	82	912149	42.712	ng	0.00
42) 2,4,6-Tribromophenol	15.778	330	775671	83.011	ng	0.00
45) 2-Fluorobiphenyl	12.860	172	2014655	40.156	ng	0.00
79) Terphenyl-d14	19.777	244	4101842	42.188	ng	-0.02

Target Compounds	Qvalue

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

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