

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061025\
Data File : BP024889.D
Acq On : 10 Jun 2025 11:33
Operator : RC/JU
Sample : PB168321BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB168321BL

Quant Time: Jun 10 12:25:37 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	238782	20.000	ng	0.00
21) Naphthalene-d8	10.378	136	925741	20.000	ng	0.00
39) Acenaphthene-d10	14.254	164	559666	20.000	ng	0.00
64) Phenanthrene-d10	17.054	188	1106475	20.000	ng	0.00
76) Chrysene-d12	21.477	240	1229440	20.000	ng	0.00
86) Perylene-d12	24.718	264	1444069	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.243	112	2193773	153.355	ng	0.00
7) Phenol-d6	6.814	99	2683147	141.761	ng	0.00
23) Nitrobenzene-d5	8.760	82	1661851	87.232	ng	0.00
42) 2,4,6-Tribromophenol	15.778	330	1164554	150.504	ng	0.00
45) 2-Fluorobiphenyl	12.854	172	3621535	87.171	ng	0.00
79) Terphenyl-d14	19.778	244	5873234	85.619	ng	-0.02

Target Compounds	Qvalue

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

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