

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061125\
 Data File : BF142724.D
 Acq On : 11 Jun 2025 09:53
 Operator : RC/JU
 Sample : PB168376BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB168376BL

Quant Time: Jun 11 11:01:19 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	6.892	152	75423	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	293240	20.000	ng	0.00
39) Acenaphthene-d10	9.933	164	164071	20.000	ng	0.00
64) Phenanthrene-d10	11.421	188	310217	20.000	ng	0.00
76) Chrysene-d12	14.068	240	203839	20.000	ng	0.00
86) Perylene-d12	15.562	264	148373	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	567406	128.127	ng	0.02
7) Phenol-d6	6.522	99	663304	127.273	ng	0.00
23) Nitrobenzene-d5	7.451	82	418446	78.095	ng	0.00
42) 2,4,6-Tribromophenol	10.727	330	237240	131.932	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	943613	76.409	ng	0.00
79) Terphenyl-d14	13.015	244	1059635	71.397	ng	0.00

Target Compounds	Qvalue

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

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