

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100625\
 Data File : BP025925.D
 Acq On : 06 Oct 2025 16:15
 Operator : CG/JU
 Sample : PB169989BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB169989BL

Quant Time: Oct 06 16:38:38 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 11 16:45:04 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.737	152	166902	20.000	ng	-0.02
21) Naphthalene-d8	10.501	136	621599	20.000	ng	-0.02
39) Acenaphthene-d10	14.360	164	409905	20.000	ng	-0.02
64) Phenanthrene-d10	17.166	188	834381	20.000	ng	-0.02
76) Chrysene-d12	21.601	240	918841	20.000	ng	-0.04
86) Perylene-d12	24.954	264	970381	20.000	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.372	112	1391143	134.491	ng	-0.01
7) Phenol-d6	6.943	99	1732470	126.009	ng	0.00
23) Nitrobenzene-d5	8.884	82	1102998	78.273	ng	-0.02
42) 2,4,6-Tribromophenol	15.878	330	689611	134.881	ng	-0.03
45) 2-Fluorobiphenyl	12.966	172	2484362	80.700	ng	-0.02
79) Terphenyl-d14	19.877	244	4166754	81.580	ng	-0.04

Target Compounds	Qvalue

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

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