

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100725\
 Data File : BP025941.D
 Acq On : 07 Oct 2025 13:01
 Operator : CG/JU
 Sample : PB169995BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB169995BL

Quant Time: Oct 07 13:25:15 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 06 19:16:54 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.737	152	161488	20.000	ng	-0.02
21) Naphthalene-d8	10.507	136	590882	20.000	ng	-0.02
39) Acenaphthene-d10	14.354	164	381445	20.000	ng	-0.02
64) Phenanthrene-d10	17.154	188	796284	20.000	ng	-0.03
76) Chrysene-d12	21.613	240	849961	20.000	ng	-0.03
86) Perylene-d12	24.971	264	886103	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.378	112	1372933	137.180	ng	0.00
7) Phenol-d6	6.943	99	1651604	124.155	ng	0.00
23) Nitrobenzene-d5	8.884	82	1077352	80.427	ng	-0.02
42) 2,4,6-Tribromophenol	15.878	330	711093	149.459	ng	-0.03
45) 2-Fluorobiphenyl	12.960	172	2375306	82.915	ng	-0.03
79) Terphenyl-d14	19.872	244	4078817	86.330	ng	-0.04

Target Compounds	Qvalue

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

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