

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080125\
 Data File : BP025290.D
 Acq On : 01 Aug 2025 12:13
 Operator : CG/JU
 Sample : PB169072BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB169072BL

Quant Time: Aug 01 12:31:17 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072125.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 29 15:28:11 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.390	152	453280	20.000	ng	-0.01
21) Naphthalene-d8	10.131	136	1653865	20.000	ng	-0.01
39) Acenaphthene-d10	14.031	164	990860	20.000	ng	-0.02
64) Phenanthrene-d10	16.860	188	1929176	20.000	ng	-0.02
76) Chrysene-d12	21.289	240	2040297	20.000	ng	-0.02
86) Perylene-d12	24.407	264	2369145	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.090	112	3324284	123.916	ng	-0.02
7) Phenol-d6	6.637	99	3873333	116.574	ng	-0.01
23) Nitrobenzene-d5	8.519	82	2427147	81.557	ng	-0.02
42) 2,4,6-Tribromophenol	15.589	330	1636743	112.048	ng	0.01
45) 2-Fluorobiphenyl	12.631	172	5876705	83.615	ng	-0.02
79) Terphenyl-d14	19.624	244	8965996	86.748	ng	0.00

Target Compounds	Qvalue

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

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