

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041025\
 Data File : BP024246.D
 Acq On : 10 Apr 2025 10:27
 Operator : RC/JU
 Sample : PB167518BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB167518BL

Quant Time: Apr 10 11:46:57 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 10 11:43:37 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.752	152	279607	20.000	ng	0.00
21) Naphthalene-d8	10.528	136	1127548	20.000	ng	0.00
39) Acenaphthene-d10	14.387	164	674843	20.000	ng	0.00
64) Phenanthrene-d10	17.181	188	1292039	20.000	ng	-0.01
76) Chrysene-d12	21.621	240	1334330	20.000	ng	-0.02
86) Perylene-d12	24.986	264	1503669	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	2662955	152.029	ng	0.01
7) Phenol-d6	6.934	99	3335740	142.409	ng	0.00
23) Nitrobenzene-d5	8.899	82	1951169	97.634	ng	0.00
42) 2,4,6-Tribromophenol	15.898	330	1351320	158.969	ng	0.00
45) 2-Fluorobiphenyl	12.998	172	4408592	96.327	ng	0.00
79) Terphenyl-d14	19.898	244	6462663	99.978	ng	-0.02

Target Compounds	Qvalue

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

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