

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092024\
 Data File : BP021960.D
 Acq On : 20 Sep 2024 09:47
 Operator : RC/JU
 Sample : PB163109BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB163109BL

Quant Time: Sep 20 11:00:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 20 01:22:00 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	93050	20.000	ng	-0.01
21) Naphthalene-d8	10.469	136	324846	20.000	ng	0.00
39) Acenaphthene-d10	14.322	164	211701	20.000	ng	0.00
64) Phenanthrene-d10	17.122	188	479935	20.000	ng	-0.02
76) Chrysene-d12	21.551	240	520357	20.000	ng	-0.01
86) Perylene-d12	24.833	264	595273	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.328	112	600629	121.305	ng	0.00
7) Phenol-d6	6.887	99	732022	113.884	ng	-0.01
23) Nitrobenzene-d5	8.846	82	495987	80.790	ng	-0.01
42) 2,4,6-Tribromophenol	15.834	330	434135	110.616	ng	-0.02
45) 2-Fluorobiphenyl	12.934	172	1221445	84.224	ng	0.00
79) Terphenyl-d14	19.851	244	2435019	87.904	ng	0.00

Target Compounds	Qvalue

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

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