

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092024\
 Data File : BP021961.D
 Acq On : 20 Sep 2024 10:28
 Operator : RC/JU
 Sample : PB163110BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB163110BL

Quant Time: Sep 20 11:01:18 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	7.716	152	61243	20.000	ng	0.00
21) Naphthalene-d8	10.469	136	204152	20.000	ng	0.00
39) Acenaphthene-d10	14.322	164	136973	20.000	ng	0.00
64) Phenanthrene-d10	17.122	188	307906	20.000	ng	-0.02
76) Chrysene-d12	21.551	240	341835	20.000	ng	-0.01
86) Perylene-d12	24.833	264	416585	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.334	112	530211	162.698	ng	0.00
7) Phenol-d6	6.899	99	626213	148.020	ng	0.00
23) Nitrobenzene-d5	8.852	82	422064	109.393	ng	0.00
42) 2,4,6-Tribromophenol	15.834	330	390775	153.889	ng	-0.02
45) 2-Fluorobiphenyl	12.934	172	1038212	110.646	ng	0.00
79) Terphenyl-d14	19.851	244	2118525	116.419	ng	0.00

Target Compounds	Qvalue

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

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