

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110824\
 Data File : BP022926.D
 Acq On : 08 Nov 2024 09:15
 Operator : RC/JU
 Sample : PB164366BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB164366BL

Quant Time: Nov 08 10:04:58 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 23:36:11 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	71448	20.000	ng	0.00
21) Naphthalene-d8	10.352	136	250478	20.000	ng	0.00
39) Acenaphthene-d10	14.210	164	189143	20.000	ng	0.00
64) Phenanthrene-d10	17.022	188	458140	20.000	ng	0.00
76) Chrysene-d12	21.451	240	568592	20.000	ng	-0.01
86) Perylene-d12	24.686	264	677061	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.234	112	545986	145.023	ng	0.00
7) Phenol-d6	6.793	99	691909	131.784	ng	0.00
23) Nitrobenzene-d5	8.740	82	481814	90.877	ng	0.00
42) 2,4,6-Tribromophenol	15.734	330	477237	130.173	ng	-0.01
45) 2-Fluorobiphenyl	12.810	172	1231267	93.015	ng	0.00
79) Terphenyl-d14	19.751	244	3041329	100.453	ng	0.00

Target Compounds	Qvalue

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

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