

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100422\
 Data File : BM036826.D
 Acq On : 04 Oct 2022 11:09
 Operator : CG/JU
 Sample : PB147982BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB147982BL

Quant Time: Oct 04 22:04:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM100322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 04 03:17:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	368073	20.000	ng	0.00
21) Naphthalene-d8	10.781	136	1573655	20.000	ng	0.00
39) Acenaphthene-d10	14.598	164	932992	20.000	ng	0.00
64) Phenanthrene-d10	17.339	188	1819410	20.000	ng	0.00
76) Chrysene-d12	21.509	240	1193161	20.000	ng	0.00
86) Perylene-d12	23.921	264	776931	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	3731205	179.906	ng	0.00
7) Phenol-d6	7.140	99	4783820	164.210	ng	0.00
23) Nitrobenzene-d5	9.140	82	2773530	94.271	ng	0.00
42) 2,4,6-Tribromophenol	16.086	330	1287196	175.908	ng	0.00
45) 2-Fluorobiphenyl	13.233	172	6083398	100.538	ng	0.00
79) Terphenyl-d14	19.956	244	7617606	127.203	ng	0.00

Target Compounds	Qvalue

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

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