

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111422\
 Data File : BP012647.D
 Acq On : 14 Nov 2022 11:12
 Operator : CG/JU
 Sample : PB148922BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB148922BL

Quant Time: Nov 14 16:57:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 14 04:00:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	293681	20.000	ng	0.00
21) Naphthalene-d8	10.504	136	1300149	20.000	ng	0.00
39) Acenaphthene-d10	14.369	164	858617	20.000	ng	0.00
64) Phenanthrene-d10	17.133	188	1852577	20.000	ng	0.00
76) Chrysene-d12	21.233	240	1906814	20.000	ng	0.00
86) Perylene-d12	23.574	264	1891169	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.293	112	2226622	141.178	ng	0.00
7) Phenol-d6	6.899	99	2901839	132.552	ng	0.00
23) Nitrobenzene-d5	8.881	82	1779984	75.954	ng	0.00
42) 2,4,6-Tribromophenol	15.875	330	1493693	127.434	ng	0.00
45) 2-Fluorobiphenyl	12.987	172	4214968	74.523	ng	0.00
79) Terphenyl-d14	19.710	244	6907916	73.394	ng	0.00

Target Compounds	Qvalue

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

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