

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111222\
 Data File : BP012637.D
 Acq On : 13 Nov 2022 01:22
 Operator : CG/JU
 Sample : N5530-07
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-10

Quant Time: Nov 14 03:25:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Nov 13 22:44:57 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.711	152	291429	20.000	ng	0.00
21) Naphthalene-d8	10.505	136	1245509	20.000	ng	-0.01
39) Acenaphthene-d10	14.369	164	833018	20.000	ng	-0.01
64) Phenanthrene-d10	17.134	188	1808447	20.000	ng	-0.01
76) Chrysene-d12	21.233	240	1867008	20.000	ng	-0.01
86) Perylene-d12	23.574	264	1770908	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.299	112	2239932	143.119	ng	0.00
7) Phenol-d6	6.899	99	3020669	139.046	ng	-0.01
23) Nitrobenzene-d5	8.881	82	1786547	79.578	ng	-0.01
42) 2,4,6-Tribromophenol	15.869	330	1627195	143.090	ng	-0.02
45) 2-Fluorobiphenyl	12.987	172	4348367	79.244	ng	-0.01
79) Terphenyl-d14	19.704	244	7211836	78.256	ng	-0.01

Target Compounds	Qvalue

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

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