

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111222\
 Data File : BP012640.D
 Acq On : 13 Nov 2022 03:05
 Operator : CG/JU
 Sample : N5531-08
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-33

Quant Time: Nov 14 03:25:59 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	7.710	152	308603	20.000	ng	0.00
21) Naphthalene-d8	10.504	136	1315034	20.000	ng	-0.01
39) Acenaphthene-d10	14.369	164	864811	20.000	ng	-0.01
64) Phenanthrene-d10	17.128	188	1852924	20.000	ng	-0.02
76) Chrysene-d12	21.233	240	1792609	20.000	ng	-0.01
86) Perylene-d12	23.574	264	1621565	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.299	112	2736947	165.143	ng	0.00
7) Phenol-d6	6.899	99	3636251	158.068	ng	-0.01
23) Nitrobenzene-d5	8.881	82	2203134	92.946	ng	-0.01
42) 2,4,6-Tribromophenol	15.869	330	1834181	155.362	ng	-0.02
45) 2-Fluorobiphenyl	12.987	172	4949061	86.875	ng	-0.01
79) Terphenyl-d14	19.704	244	7678888	86.783	ng	-0.01

Target Compounds Qvalue

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

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