

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112322\
 Data File : BP012714.D
 Acq On : 23 Nov 2022 17:14
 Operator : CG/JU
 Sample : N5749-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-82

Quant Time: Nov 24 02:31:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 23 00:38:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	354236	20.000	ng	0.00
21) Naphthalene-d8	10.846	136	1319108	20.000	ng	0.00
39) Acenaphthene-d10	14.663	164	744590	20.000	ng	0.00
64) Phenanthrene-d10	17.422	188	1546641	20.000	ng	0.00
76) Chrysene-d12	21.504	240	1690780	20.000	ng	0.00
86) Perylene-d12	24.027	264	1780228	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	2982436	182.400	ng	0.00
7) Phenol-d6	7.181	99	3380685	144.168	ng	0.00
23) Nitrobenzene-d5	9.199	82	2135590	90.291	ng	0.00
42) 2,4,6-Tribromophenol	16.157	330	1452272	148.894	ng	0.00
45) 2-Fluorobiphenyl	13.293	172	4846373	92.728	ng	0.00
79) Terphenyl-d14	19.969	244	8658518	108.840	ng	0.00
Target Compounds						
32) Benzoic acid	10.069	122	127164	22.661	ng	Qvalue 97
74) Di-n-butylphthalate	18.369	149	13131	5.845	ng	98

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

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