

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030922\
 Data File : BP009329.D
 Acq On : 09 Mar 2022 18:09
 Operator : CG/JU
 Sample : N1830-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 11986

Quant Time: Mar 10 04:10:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 01:55:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	332846	20.000	ng	0.00
21) Naphthalene-d8	10.610	136	1488087	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	1048008	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	2517593	20.000	ng	0.00
76) Chrysene-d12	21.322	240	2993628	20.000	ng	0.00
86) Perylene-d12	23.727	264	2898922	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.411	112	2982637	130.802	ng	0.00
7) Phenol-d6	6.993	99	3724722	128.539	ng	0.00
23) Nitrobenzene-d5	8.969	82	2664138	98.608	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	2372245	187.459	ng	0.00
45) 2-Fluorobiphenyl	13.081	172	6642958	87.170	ng	0.00
79) Terphenyl-d14	19.792	244	12962475	81.713	ng	0.00

Target Compounds	Qvalue

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

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