

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032922\
 Data File : BP009597.D
 Acq On : 29 Mar 2022 16:35
 Operator : CG/JU
 Sample : PB143709BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB143709BL

Quant Time: Mar 30 04:09:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 30 04:00:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	252750	20.000	ng	0.00
21) Naphthalene-d8	10.557	136	951618	20.000	ng	0.00
39) Acenaphthene-d10	14.410	164	553244	20.000	ng	0.00
64) Phenanthrene-d10	17.174	188	1114403	20.000	ng	0.00
76) Chrysene-d12	21.292	240	1127771	20.000	ng	0.00
86) Perylene-d12	23.686	264	1237941	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	2162622	149.388	ng	0.00
7) Phenol-d6	6.951	99	2617617	133.706	ng	0.00
23) Nitrobenzene-d5	8.922	82	1651167	92.205	ng	0.00
42) 2,4,6-Tribromophenol	15.910	330	1016417	111.333	ng	0.00
45) 2-Fluorobiphenyl	13.027	172	3746654	93.882	ng	0.00
79) Terphenyl-d14	19.757	244	5510763	87.706	ng	0.00

Target Compounds	Qvalue

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

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