

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
 Data File : BP009421.D
 Acq On : 16 Mar 2022 13:51
 Operator : CG/JU
 Sample : PB142486BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB142486BL

Quant Time: Mar 16 15:33:00 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	7.811	152	306526	20.000	ng	-0.01
21) Naphthalene-d8	10.605	136	1169187	20.000	ng	-0.01
39) Acenaphthene-d10	14.451	164	746518	20.000	ng	-0.01
64) Phenanthrene-d10	17.210	188	1547325	20.000	ng	-0.01
76) Chrysene-d12	21.322	240	1393621	20.000	ng	0.00
86) Perylene-d12	23.727	264	1429168	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.417	112	2375882	113.140	ng	0.00
7) Phenol-d6	7.005	99	2930611	109.818	ng	0.02
23) Nitrobenzene-d5	8.963	82	1882862	88.699	ng	-0.01
42) 2,4,6-Tribromophenol	15.951	330	1522961	168.951	ng	0.00
45) 2-Fluorobiphenyl	13.069	172	4589734	84.550	ng	-0.02
79) Terphenyl-d14	19.786	244	6827998	92.459	ng	-0.01

Target Compounds	Qvalue

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

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