

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
 Data File : BP019842.D
 Acq On : 23 Feb 2024 14:25
 Operator : MA/JU
 Sample : P1521-04DL 50X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CG-MW-17-CH3DL

Quant Time: Feb 23 14:54:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 23 13:10:49 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	61105	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	216362	20.000	ng	0.00
39) Acenaphthene-d10	14.545	164	149960	20.000	ng	0.00
64) Phenanthrene-d10	17.310	188	324183	20.000	ng	0.00
76) Chrysene-d12	21.416	240	360774	20.000	ng	-0.01
86) Perylene-d12	23.851	264	493740	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	4890	1.290	ng	0.00
7) Phenol-d6	7.087	99	3642	0.713	ng	0.01
23) Nitrobenzene-d5	9.075	82	8357	1.674	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	6798	3.105	ng	0.00
45) 2-Fluorobiphenyl	13.175	172	18157	1.497	ng	0.00
79) Terphenyl-d14	19.880	244	33214	1.514	ng	-0.01
Target Compounds						
20) 3+4-Methylphenols	8.675	107	164444	35.132	ng	Qvalue 89

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

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