

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021924\
 Data File : BP019769.D
 Acq On : 19 Feb 2024 18:36
 Operator : MA/JU
 Sample : P1505-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 WC-1D

Quant Time: Feb 19 23:37:03 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 19 23:35:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	52167	20.000	ng	0.00
21) Naphthalene-d8	10.698	136	193191	20.000	ng	0.00
39) Acenaphthene-d10	14.534	164	134083	20.000	ng	0.00
64) Phenanthrene-d10	17.292	188	339988	20.000	ng	0.00
76) Chrysene-d12	21.380	240	374409	20.000	ng	-0.01
86) Perylene-d12	23.804	264	401376	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	415185	128.289	ng	0.00
7) Phenol-d6	7.069	99	481850	110.421	ng	0.00
23) Nitrobenzene-d5	9.069	82	401256	89.999	ng	0.00
42) 2,4,6-Tribromophenol	16.028	330	379816	194.011	ng	0.00
45) 2-Fluorobiphenyl	13.163	172	947938	87.399	ng	0.00
79) Terphenyl-d14	19.857	244	2048833	89.992	ng	0.00
Target Compounds						
88) Benzo(b)fluoranthene	23.063	252	50823m	2.009	ng	Qvalue
89) Benzo(k)fluoranthene	23.110	252	58752	2.431	ng	98

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

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