

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020924\
 Data File : BP019653.D
 Acq On : 10 Feb 2024 02:56
 Operator : MA/JU
 Sample : P1378-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-4

Quant Time: Feb 10 03:30:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 01 00:51:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	98604	20.000	ng	-0.01
21) Naphthalene-d8	10.704	136	349056	20.000	ng	-0.01
39) Acenaphthene-d10	14.539	164	191288	20.000	ng	-0.01
64) Phenanthrene-d10	17.298	188	388038	20.000	ng	0.00
76) Chrysene-d12	21.386	240	457317	20.000	ng	0.00
86) Perylene-d12	23.809	264	553514	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	502488	82.144	ng	0.00
7) Phenol-d6	7.081	99	730793	88.601	ng	0.00
23) Nitrobenzene-d5	9.075	82	481878	59.820	ng	-0.01
42) 2,4,6-Tribromophenol	16.033	330	337063	120.684	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	1091558	70.544	ng	-0.01
79) Terphenyl-d14	19.862	244	1773294	63.769	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.304	202	66231	2.655	ng	99
78) Pyrene	19.657	202	90726	2.811	ng	99
81) Benzo(a)anthracene	21.374	228	66024	2.051	ng	95
83) Chrysene	21.421	228	68018m	2.225	ng	
88) Benzo(b)fluoranthene	23.068	252	91121m	2.612	ng	
90) Benzo(a)pyrene	23.698	252	73992	2.384	ng	93

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

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