

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082224\
 Data File : BP021619.D
 Acq On : 22 Aug 2024 18:41
 Operator : RC/JU
 Sample : P3518-17
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OU4-TS-09-080724

Quant Time: Aug 23 00:53:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 11:12:31 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	352434	20.000	ng	0.00
21) Naphthalene-d8	10.587	136	1405074	20.000	ng	0.00
39) Acenaphthene-d10	14.445	164	872533	20.000	ng	0.00
64) Phenanthrene-d10	17.257	188	1872743	20.000	ng	0.00
76) Chrysene-d12	21.716	240	1824677	20.000	ng	0.00
86) Perylene-d12	25.145	264	2134055	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.399	112	2236247	94.217	ng	0.00
7) Phenol-d6	6.969	99	3117879	90.049	ng	0.00
23) Nitrobenzene-d5	8.946	82	1851453	68.013	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	1016626	104.689	ng	0.00
45) 2-Fluorobiphenyl	13.057	172	2841645	49.704	ng	0.00
79) Terphenyl-d14	19.980	244	4081360	43.454	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.392	202	357135	3.232	ng	99
78) Pyrene	19.763	202	347831	2.916	ng	100
88) Benzo(b)fluoranthene	24.063	252	311178	2.539	ng	# 93

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

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