

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022924\
 Data File : BP019946.D
 Acq On : 29 Feb 2024 20:13
 Operator : MA/JU
 Sample : P1639-31
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP-8

Quant Time: Mar 01 03:44:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP022624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 29 12:18:35 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	70223	20.000	ng	-0.02
21) Naphthalene-d8	10.693	136	281614	20.000	ng	-0.02
39) Acenaphthene-d10	14.540	164	216054	20.000	ng	-0.02
64) Phenanthrene-d10	17.298	188	563322	20.000	ng	-0.02
76) Chrysene-d12	21.404	240	594935	20.000	ng	-0.02
86) Perylene-d12	23.833	264	603316	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	360331	84.223	ng	-0.02
7) Phenol-d6	7.069	99	508534	79.835	ng	-0.02
23) Nitrobenzene-d5	9.063	82	337872	48.436	ng	-0.02
42) 2,4,6-Tribromophenol	16.039	330	376552	101.645	ng	-0.02
45) 2-Fluorobiphenyl	13.163	172	818851	51.207	ng	-0.02
79) Terphenyl-d14	19.875	244	2073866	54.942	ng	-0.02
Target Compounds						
75) Fluoranthene	19.316	202	158975	4.074	ng	99
78) Pyrene	19.669	202	134872	3.300	ng	99
88) Benzo(b)fluoranthene	23.092	252	96257	2.602	ng	# 92

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

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