

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022324\
 Data File : BP019849.D
 Acq On : 23 Feb 2024 18:21
 Operator : MA/JU
 Sample : P1537-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 13

Quant Time: Feb 24 00:02:27 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	7.905	152	79484	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	340328	20.000	ng	0.00
39) Acenaphthene-d10	14.545	164	195175	20.000	ng	0.00
64) Phenanthrene-d10	17.310	188	440049	20.000	ng	0.00
76) Chrysene-d12	21.416	240	437048	20.000	ng	-0.01
86) Perylene-d12	23.851	264	515803	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	468109	94.932	ng	0.00
7) Phenol-d6	7.081	99	606425	91.208	ng	0.00
23) Nitrobenzene-d5	9.075	82	516328	65.740	ng	0.00
42) 2,4,6-Tribromophenol	16.045	330	355206	124.647	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	966870	61.241	ng	0.00
79) Terphenyl-d14	19.880	244	1721390	64.773	ng	-0.01
Target Compounds						
						Qvalue
75) Fluoranthene	19.322	202	136715	4.833	ng	100
78) Pyrene	19.675	202	123628	4.008	ng	99
81) Benzo(a)anthracene	21.398	228	69375	2.255	ng	97
83) Chrysene	21.451	228	78346	2.682	ng	95
88) Benzo(b)fluoranthene	23.110	252	117994	3.630	ng	# 92
90) Benzo(a)pyrene	23.739	252	80015	2.766	ng	# 89
92) Benzo(g,h,i)perylene	27.180	276	69649	2.114	ng	95

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

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