

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081624\
 Data File : BP021540.D
 Acq On : 16 Aug 2024 15:12
 Operator : RC/JU
 Sample : P3518-05RE
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OU4-TS-03-080724RE

Quant Time: Aug 16 15:41:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 14 00:00:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	543985	20.000	ng	0.01
21) Naphthalene-d8	10.610	136	2309246	20.000	ng	0.01
39) Acenaphthene-d10	14.463	164	1337933	20.000	ng	0.00
64) Phenanthrene-d10	17.269	188	2187436	20.000	ng	0.02
76) Chrysene-d12	21.715	240	1743108	20.000	ng	0.02
86) Perylene-d12	25.133	264	2225455	20.000	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	2996653	81.797	ng	0.02
7) Phenol-d6	6.981	99	4380194	81.961	ng	0.02
23) Nitrobenzene-d5	8.963	82	2544197	56.867	ng	0.01
42) 2,4,6-Tribromophenol	15.975	330	1054660	70.827	ng	0.00
45) 2-Fluorobiphenyl	13.075	172	3718495	42.416	ng	0.01
79) Terphenyl-d14	19.986	244	3375764	37.624	ng	0.01
Target Compounds						
						Qvalue
75) Fluoranthene	19.398	202	386670	2.996	ng	99
78) Pyrene	19.768	202	356920	3.133	ng	98
83) Chrysene	21.768	228	235780	2.282	ng	# 94
88) Benzo(b)fluoranthene	24.051	252	381660	2.986	ng	# 92
90) Benzo(a)pyrene	24.974	252	291358	2.627	ng	# 92

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

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