

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081624\
 Data File : BP021534.D
 Acq On : 16 Aug 2024 10:59
 Operator : RC/JU
 Sample : P3518-01RE
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OU4-TS-01-080724RE

Quant Time: Aug 16 11:59:12 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	453420	20.000	ng	0.01
21) Naphthalene-d8	10.604	136	1930500	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	1213530	20.000	ng	0.01
64) Phenanthrene-d10	17.269	188	2352082	20.000	ng	0.02
76) Chrysene-d12	21.721	240	1990799	20.000	ng	0.02
86) Perylene-d12	25.139	264	2355058	20.000	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.411	112	1846623	60.473	ng	0.02
7) Phenol-d6	6.975	99	2612195	58.641	ng	0.01
23) Nitrobenzene-d5	8.963	82	1638951	43.820	ng	0.01
42) 2,4,6-Tribromophenol	15.981	330	796540	58.976	ng	0.01
45) 2-Fluorobiphenyl	13.075	172	2490324	31.319	ng	0.01
79) Terphenyl-d14	19.992	244	2947460	28.763	ng	0.02
Target Compounds						
						Qvalue
75) Fluoranthene	19.392	202	358610	2.584	ng	100
78) Pyrene	19.769	202	343687	2.641	ng	100

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

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