

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081624\
 Data File : BP021541.D
 Acq On : 16 Aug 2024 15:54
 Operator : RC/JU
 Sample : P3518-07RE
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OU4-TS-04-080724RE

Quant Time: Aug 16 16:21:44 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	467262	20.000	ng	0.01
21) Naphthalene-d8	10.610	136	1831532	20.000	ng	0.01
39) Acenaphthene-d10	14.463	164	983158	20.000	ng	0.00
64) Phenanthrene-d10	17.269	188	1814668	20.000	ng	0.02
76) Chrysene-d12	21.716	240	1720162	20.000	ng	0.02
86) Perylene-d12	25.145	264	2107755	20.000	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.411	112	2142603	68.087	ng	0.02
7) Phenol-d6	6.981	99	3020891	65.807	ng	0.02
23) Nitrobenzene-d5	8.963	82	1732640	48.828	ng	0.01
42) 2,4,6-Tribromophenol	15.975	330	673344	61.537	ng	0.00
45) 2-Fluorobiphenyl	13.069	172	2273211	35.287	ng	0.00
79) Terphenyl-d14	19.980	244	2751973	31.080	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.386	202	274944	2.568	ng	100
78) Pyrene	19.763	202	247156	2.198	ng	100

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

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