

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030922\
 Data File : BP009334.D
 Acq On : 09 Mar 2022 20:58
 Operator : CG/JU
 Sample : N1762-01DL2 50X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 282DL2

Quant Time: Mar 10 04:12:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 01:55:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	311600	20.000	ng	0.00
21) Naphthalene-d8	10.610	136	1442175	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	1038038	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	2471029	20.000	ng	0.00
76) Chrysene-d12	21.321	240	3225805	20.000	ng	0.00
86) Perylene-d12	23.727	264	3737459	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	30045	1.407	ng	0.00
7) Phenol-d6	6.987	99	41512	1.530	ng	0.00
23) Nitrobenzene-d5	8.963	82	26713	1.020	ng	-0.01
42) 2,4,6-Tribromophenol	15.945	330	21030	1.678	ng	-0.01
45) 2-Fluorobiphenyl	13.081	172	71024	0.941	ng	0.00
79) Terphenyl-d14	19.786	244	165218	0.967	ng	-0.01
Target Compounds						
						Qvalue
31) Naphthalene	10.657	128	1270637	15.910	ng	99
37) 2-Methylnaphthalene	12.275	142	717015	12.636	ng	99
38) 1-Methylnaphthalene	12.492	142	372422	7.126	ng	99
46) 1,1'-Biphenyl	13.287	154	244809	2.889	ng	98
52) Acenaphthene	14.516	154	1185050	18.642	ng	99
55) Dibenzofuran	14.851	168	1278855	13.281	ng	99
58) Fluorene	15.504	166	1249153	16.565	ng	98
71) Phenanthrene	17.257	178	6287223	44.273	ng	99
72) Anthracene	17.345	178	417190	2.951	ng	97
75) Fluoranthene	19.233	202	3959124	24.154	ng	99
78) Pyrene	19.586	202	2797768	12.846	ng	99
81) Benzo(a)anthracene	21.304	228	720044	3.293	ng	98
83) Chrysene	21.357	228	656694	3.093	ng	98

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

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