

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE040821\
 Data File : BE101261.D
 Acq On : 9 Apr 2021 4:31
 Operator : CG/JU
 Sample : M1967-13
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 EB 2021-04-06 SED

Quant Time: Apr 09 05:14:48 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-SIM-BE040621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 07 10:22:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.838	152	3108	0.40	ng	0.00
7) Naphthalene-d8	10.615	136	14262	0.40	ng	0.00
13) Acenaphthene-d10	14.443	164	10506	0.40	ng	0.00
19) Phenanthrene-d10	17.195	188	31138	0.40	ng	#-0.01
29) Chrysene-d12	21.354	240	25768	0.40	ng	0.00
36) Perylene-d12	23.838	264	25679	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.431	112	1509	0.21	ng	0.00
5) Phenol-d6	6.997	99	1370	0.12	ng	0.00
8) Nitrobenzene-d5	8.979	82	4345	0.41	ng	0.00
11) 2-Methylnaphthalene-d10	12.196	152	7504	0.23	ng	0.00
14) 2,4,6-Tribromophenol	15.955	330	867	0.31	ng	0.00
15) 2-Fluorobiphenyl	13.075	172	11795	0.31	ng	0.00
27) Fluoranthene-d10	19.215	212	142793	0.26	ng	0.00
31) Terphenyl-d14	19.812	244	36604	0.53	ng	0.00
Target Compounds						Qvalue
32) Benzo(a)anthracene	21.342	228	1542	0.02	ng	# 84
34) Bis(2-ethylhexyl)phtha...	21.281	149	22598	0.43	ng	99
35) Indeno(1,2,3-cd)pyrene	26.445	276	1538	0.02	ng	# 87

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

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