

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012622\
 Data File : BM033902.D
 Acq On : 27 Jan 2022 14:55
 Operator : CG/JU
 Sample : PB142306BS
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS306

Quant Time: Jan 28 01:50:51 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM012422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 24 15:20:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.939	152	1852	0.400	ng/ul	0.00
4) Naphthalene-d8	10.751	136	6300	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.576	164	2805	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	5594	0.400	ng/ul	0.00
17) Chrysene-d12	21.477	240	4808	0.400	ng/ul #	0.00
23) Perylene-d12	23.873	264	4699	0.400	ng/ul #-	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.292	96	1293	0.668	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.344	152	2720	0.313	ng/ul	0.00
18) Fluoranthene-d10	19.334	212	5199	0.351	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.330	88	3227	1.592	ng/ul#	78
5) Naphthalene	10.801	128	5670	0.327	ng/ul#	90
7) 2-Methylnaphthalene	12.416	142	3274	0.291	ng/ul	98
8) 1-Methylnaphthalene	12.631	142	3206	0.291	ng/ul	99
10) Acenaphthylene	14.299	152	4385	0.350	ng/ul	93
11) Acenaphthene	14.640	153	3210	0.320	ng/ul	97
12) Fluorene	15.630	166	3610	0.315	ng/ul	97
14) Pentachlorophenol	16.972	266	1354	0.527	ng/ul	98
15) Phenanthrene	17.358	178	5856	0.329	ng/ul	98
16) Anthracene	17.451	178	5234	0.314	ng/ul	96
19) Fluoranthene	19.365	202	7332	0.335	ng/ul	99
20) Pyrene	19.723	202	7684	0.342	ng/ul#	94
21) Benzo(a)anthracene	21.463	228	6367	0.324	ng/ul	99
22) Chrysene	21.514	228	7640	0.367	ng/ul	98
24) Benzo(b)fluoranthene	23.141	252	7394	0.350	ng/ul	94
25) Benzo(k)fluoranthene	23.187	252	7886	0.373	ng/ul#	94
26) Benzo(a)pyrene	23.769	252	7102	0.388	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	26.385	276	8486	0.349	ng/ul#	87
28) Dibenzo(a,h)anthracene	26.410	278	6707	0.349	ng/ul	97
29) Benzo(g,h,i)perylene	27.141	276	7766	0.376	ng/ul	96

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

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