

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042522\
 Data File : BM034803.D
 Acq On : 25 Apr 2022 17:39
 Operator : CG/JU
 Sample : PB144306BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS306

Quant Time: Apr 26 03:05:41 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM042222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 22 16:54:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.942	152	6028	0.400	ng/ul	0.00
4) Naphthalene-d8	10.743	136	15112	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.571	164	7378	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.310	188	13231	0.400	ng/ul #	0.00
17) Chrysene-d12	21.487	240	9007	0.400	ng/ul #	0.00
23) Perylene-d12	23.881	264	9060	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.386	96	4485	0.673	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.332	152	7643	0.360	ng/ul	0.00
18) Fluoranthene-d10	19.336	212	11427	0.402	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.420	88	11107	1.750	ng/ul#	81
5) Naphthalene	10.792	128	15367	0.356	ng/ul#	87
7) 2-Methylnaphthalene	12.404	142	9376	0.327	ng/ul	99
8) 1-Methylnaphthalene	12.624	142	9218	0.327	ng/ul	100
10) Acenaphthylene	14.289	152	12851	0.399	ng/ul#	90
11) Acenaphthene	14.631	153	9532	0.349	ng/ul	97
12) Fluorene	15.617	166	10329	0.336	ng/ul#	97
14) Pentachlorophenol	16.964	266	2826	0.720	ng/ul	98
15) Phenanthrene	17.352	178	15689	0.359	ng/ul	95
16) Anthracene	17.445	178	13847	0.355	ng/ul#	92
19) Fluoranthene	19.364	202	16586	0.384	ng/ul	97
20) Pyrene	19.727	202	16474	0.379	ng/ul	99
21) Benzo(a)anthracene	21.473	228	14058	0.401	ng/ul	98
22) Chrysene	21.525	228	14961	0.400	ng/ul	98
24) Benzo(b)fluoranthene	23.153	252	15764	0.381	ng/ul	91
25) Benzo(k)fluoranthene	23.200	252	15092	0.365	ng/ul#	90
26) Benzo(a)pyrene	23.776	252	14241	0.422	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.362	276	18213	0.384	ng/ul#	83
28) Dibenzo(a,h)anthracene	26.379	278	14639	0.380	ng/ul#	91
29) Benzo(g,h,i)perylene	27.117	276	15890	0.388	ng/ul	93

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

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