

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042222\
 Data File : BM034790.D
 Acq On : 23 Apr 2022 04:36
 Operator : CG/JU
 Sample : PB144176BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS176

Quant Time: Apr 23 05:21:37 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	7.943	152	7457	0.400	ng/ul	0.00
4) Naphthalene-d8	10.743	136	19374	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.575	164	9732	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.313	188	17179	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.485	240	12107	0.400	ng/ul	# 0.00
23) Perylene-d12	23.879	264	12884	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.386	96	4904	0.595	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.332	152	9943	0.365	ng/ul	0.00
18) Fluoranthene-d10	19.336	212	14858	0.389	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.420	88	11797	1.503	ng/ul#	82
5) Naphthalene	10.793	128	19323	0.349	ng/ul#	88
7) 2-Methylnaphthalene	12.404	142	11982	0.326	ng/ul	100
8) 1-Methylnaphthalene	12.624	142	11756	0.325	ng/ul	99
10) Acenaphthylene	14.293	152	16916	0.398	ng/ul#	90
11) Acenaphthene	14.635	153	12103	0.336	ng/ul	95
12) Fluorene	15.621	166	12770	0.315	ng/ul#	97
14) Pentachlorophenol	16.963	266	3737	0.733	ng/ul	97
15) Phenanthrene	17.351	178	19067	0.336	ng/ul	95
16) Anthracene	17.444	178	17199	0.340	ng/ul#	92
19) Fluoranthene	19.368	202	20471	0.352	ng/ul	96
20) Pyrene	19.726	202	20542	0.352	ng/ul	99
21) Benzo(a)anthracene	21.471	228	18561	0.394	ng/ul	98
22) Chrysene	21.523	228	18848	0.375	ng/ul	98
24) Benzo(b)fluoranthene	23.151	252	20183	0.343	ng/ul	91
25) Benzo(k)fluoranthene	23.198	252	19541	0.332	ng/ul#	90
26) Benzo(a)pyrene	23.771	252	19522	0.407	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	26.360	276	23899	0.355	ng/ul#	83
28) Dibenzo(a,h)anthracene	26.380	278	19321	0.352	ng/ul#	90
29) Benzo(g,h,i)perylene	27.118	276	20608	0.354	ng/ul	94

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

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