

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070821\
 Data File : BM030917.D
 Acq On : 09 Jul 2021 12:16
 Operator : CG/JU
 Sample : PB137404BS
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS404

Quant Time: Jul 09 12:48:12 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 08 10:44:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.767	152	1872	0.400	ng/ul	0.00
4) Naphthalene-d8	10.559	136	6995	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.406	164	3206	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.155	188	5576	0.400	ng/ul	0.00
17) Chrysene-d12	21.319	240	4726	0.400	ng/ul	0.00
23) Perylene-d12	23.566	264	4076	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.265	96	1583	0.776	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.164	152	3606	0.328	ng/ul	0.00
18) Fluoranthene-d10	19.169	212	5906	0.436	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.296	88	3892	1.847	ng/ul#	83
5) Naphthalene	10.613	128	6812	0.342	ng/ul	97
7) 2-Methylnaphthalene	12.241	142	4150	0.309	ng/ul	97
8) 1-Methylnaphthalene	12.452	142	3776	0.289	ng/ul	99
10) Acenaphthylene	14.129	152	4841	0.338	ng/ul	94
11) Acenaphthene	14.470	153	4224	0.364	ng/ul	98
12) Fluorene	15.471	166	4267	0.334	ng/ul	96
14) Pentachlorophenol	16.825	266	1201	0.660	ng/ul	96
15) Phenanthrene	17.193	178	6312	0.337	ng/ul	96
16) Anthracene	17.297	178	5538	0.337	ng/ul	95
19) Fluoranthene	19.199	202	7642	0.366	ng/ul	98
20) Pyrene	19.558	202	8134	0.388	ng/ul	98
21) Benzo(a)anthracene	21.309	228	5888	0.313	ng/ul	99
22) Chrysene	21.355	228	7968	0.378	ng/ul	98
24) Benzo(b)fluoranthene	22.890	252	6538	0.362	ng/ul	87
25) Benzo(k)fluoranthene	22.936	252	8267	0.444	ng/ul#	88
26) Benzo(a)pyrene	23.476	252	6002	0.391	ng/ul#	80
27) Indeno(1,2,3-cd)pyrene	25.865	276	7781	0.389	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.885	278	6318	0.394	ng/ul#	82
29) Benzo(g,h,i)perylene	26.544	276	7081	0.407	ng/ul	97

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

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