

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102521\
 Data File : BM032706.D
 Acq On : 25 Oct 2021 22:49
 Operator : CG/JU
 Sample : PB140193BS
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS193

Quant Time: Oct 26 01:42:59 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM101421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 25 10:44:54 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.546	152	4242	0.400	ng/ul	0.00
4) Naphthalene-d8	10.326	136	14961	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.198	164	7686	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.932	188	13970	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.109	240	9045	0.400	ng/ul	0.00
23) Perylene-d12	23.240	264	8080	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.917	96	3127	0.658	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.922	152	6022	0.302	ng/ul	0.00
18) Fluoranthene-d10	18.958	212	9683	0.340	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.951	88	8900	1.699	ng/ul#	80
5) Naphthalene	10.375	128	13386	0.295	ng/ul	99
7) 2-Methylnaphthalene	11.994	142	8487	0.286	ng/ul	99
8) 1-Methylnaphthalene	12.219	142	8158	0.278	ng/ul	99
10) Acenaphthylene	13.915	152	9051	0.276	ng/ul	100
11) Acenaphthene	14.259	153	9008	0.301	ng/ul	99
12) Fluorene	15.248	166	9691	0.287	ng/ul	99
14) Pentachlorophenol	16.606	266	2100	0.584	ng/ul	99
15) Phenanthrene	16.974	178	13982	0.298	ng/ul	100
16) Anthracene	17.064	178	11263	0.273	ng/ul	99
19) Fluoranthene	18.989	202	14496	0.324	ng/ul	99
20) Pyrene	19.351	202	14623	0.318	ng/ul	100
21) Benzo(a)anthracene	21.094	228	10132	0.287	ng/ul	99
22) Chrysene	21.145	228	12168	0.313	ng/ul	100
24) Benzo(b)fluoranthene	22.607	252	11667	0.313	ng/ul	88
25) Benzo(k)fluoranthene	22.648	252	12947	0.348	ng/ul#	88
26) Benzo(a)pyrene	23.150	252	9676	0.309	ng/ul#	82
27) Indeno(1,2,3-cd)pyrene	25.340	276	12985	0.339	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.347	278	10237	0.339	ng/ul#	92
29) Benzo(g,h,i)perylene	25.982	276	12033	0.361	ng/ul	99

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

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