

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070821\
 Data File : BM030890.D
 Acq On : 08 Jul 2021 18:23
 Operator : CG/JU
 Sample : PB137404BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS404

Quant Time: Jul 09 02:52:35 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	1843	0.400	ng/ul	0.00
4) Naphthalene-d8	10.559	136	7169	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.402	164	3585	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.148	188	5925	0.400	ng/ul	0.00
17) Chrysene-d12	21.311	240	4293	0.400	ng/ul	# 0.00
23) Perylene-d12	23.561	264	3876	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	1723	0.858	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.160	152	4551	0.404	ng/ul	0.00
18) Fluoranthene-d10	19.165	212	6443	0.523	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.296	88	4187	2.018	ng/ul#	83
5) Naphthalene	10.608	128	8110	0.397	ng/ul	98
7) 2-Methylnaphthalene	12.236	142	5199	0.378	ng/ul	98
8) 1-Methylnaphthalene	12.448	142	4762	0.356	ng/ul	99
10) Acenaphthylene	14.125	152	6175	0.386	ng/ul	96
11) Acenaphthene	14.466	153	5304	0.409	ng/ul	99
12) Fluorene	15.464	166	5423	0.379	ng/ul	98
14) Pentachlorophenol	16.815	266	988	0.511	ng/ul	96
15) Phenanthrene	17.190	178	7575	0.380	ng/ul	97
16) Anthracene	17.287	178	6764	0.388	ng/ul	97
19) Fluoranthene	19.195	202	8464	0.447	ng/ul	100
20) Pyrene	19.558	202	8728	0.458	ng/ul	99
21) Benzo(a)anthracene	21.299	228	6433	0.376	ng/ul	100
22) Chrysene	21.348	228	7991	0.417	ng/ul	97
24) Benzo(b)fluoranthene	22.883	252	7679	0.447	ng/ul	90
25) Benzo(k)fluoranthene	22.927	252	8616	0.487	ng/ul#	91
26) Benzo(a)pyrene	23.467	252	6577	0.450	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	25.846	276	8975	0.472	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.865	278	7275	0.477	ng/ul#	90
29) Benzo(g,h,i)perylene	26.532	276	7952	0.481	ng/ul	97

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

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