

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092721\
 Data File : BM032199.D
 Acq On : 27 Sep 2021 10:37
 Operator : CG/JU
 Sample : SSTDO.230
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDO.2030

Quant Time: Sep 27 11:48:18 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM092721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 27 11:46:23 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.664	152	4112	0.400	ng/ul	0.00
4) Naphthalene-d8	10.456	136	13274	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.308	164	6977	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.033	188	15405	0.400	ng/ul	0.00
17) Chrysene-d12	21.196	240	15368	0.400	ng/ul	0.00
23) Perylene-d12	23.363	264	13540	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.989	96	953	0.211	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.044	152	3611	0.176	ng/ul	0.00
18) Fluoranthene-d10	19.053	212	7731	0.172	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.027	88	882	0.200	ng/ul#	79
5) Naphthalene	10.505	128	8102	0.230	ng/ul	98
7) 2-Methylnaphthalene	12.120	142	5212	0.212	ng/ul	100
8) 1-Methylnaphthalene	12.340	142	5053	0.209	ng/ul	99
10) Acenaphthylene	14.023	152	5664	0.201	ng/ul	97
11) Acenaphthene	14.372	153	5287	0.230	ng/ul	98
12) Fluorene	15.351	166	5997	0.223	ng/ul	98
14) Pentachlorophenol	16.700	266	790	0.194	ng/ul	98
15) Phenanthrene	17.074	178	10210	0.218	ng/ul	98
16) Anthracene	17.165	178	8284	0.201	ng/ul	97
19) Fluoranthene	19.084	202	11819	0.178	ng/ul	99
20) Pyrene	19.442	202	12106	0.177	ng/ul	99
21) Benzo(a)anthracene	21.181	228	10885	0.194	ng/ul	99
22) Chrysene	21.232	228	13071	0.203	ng/ul	99
24) Benzo(b)fluoranthene	22.716	252	14012	0.225	ng/ul	87
25) Benzo(k)fluoranthene	22.757	252	13334	0.208	ng/ul#	86
26) Benzo(a)pyrene	23.266	252	10541	0.194	ng/ul#	83
27) Indeno(1,2,3-cd)pyrene	25.505	276	13923	0.188	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.507	278	11182	0.188	ng/ul	94
29) Benzo(g,h,i)perylene	26.166	276	13489	0.205	ng/ul	96

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

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