

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090321\
 Data File : BM032082.D
 Acq On : 04 Sep 2021 17:52
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4122

Quant Time: Sep 06 03:51:01 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM090321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 03 16:21:27 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.452	152	1735	0.400	ng/ul	0.00
4) Naphthalene-d8	10.208	136	5897	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.095	164	2513	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	16.855	188	4988	0.400	ng/ul	0.00
17) Chrysene-d12	21.061	240	3872	0.400	ng/ul	-0.01
23) Perylene-d12	23.175	264	4175	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.068	96	770	0.404	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.809	152	3073	0.337	ng/ul	-0.01
18) Fluoranthene-d10	18.896	212	4338	0.382	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.099	88	757	0.407	ng/ul#	81
5) Naphthalene	10.257	128	6256	0.400	ng/ul	97
7) 2-Methylnaphthalene	11.885	142	3773	0.346	ng/ul	98
8) 1-Methylnaphthalene	12.106	142	3635	0.338	ng/ul	98
10) Acenaphthylene	13.815	152	5173	0.510	ng/ul	99
11) Acenaphthene	14.159	153	3665	0.442	ng/ul#	87
12) Fluorene	15.160	166	4072	0.421	ng/ul	95
14) Pentachlorophenol	16.525	266	618	0.468	ng/ul	97
15) Phenanthrene	16.897	178	6034	0.397	ng/ul	99
16) Anthracene	16.990	178	5059	0.380	ng/ul	98
19) Fluoranthene	18.923	202	6280	0.375	ng/ul#	96
20) Pyrene	19.289	202	5968	0.346	ng/ul#	95
21) Benzo(a)anthracene	21.047	228	5478	0.387	ng/ul	99
22) Chrysene	21.098	228	6032	0.371	ng/ul	100
24) Benzo(b)fluoranthene	22.550	252	7791	0.406	ng/ul	94
25) Benzo(k)fluoranthene	22.591	252	7259	0.367	ng/ul	94
26) Benzo(a)pyrene	23.086	252	6904	0.412	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	25.242	276	9228	0.404	ng/ul#	92
28) Dibenzo(a,h)anthracene	25.257	278	7279	0.396	ng/ul	94
29) Benzo(g,h,i)perylene	25.874	276	7982	0.393	ng/ul	96

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

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