

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062521\
 Data File : BM030646.D
 Acq On : 26 Jun 2021 20:47
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4709

Quant Time: Jun 27 02:43:02 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 25 10:09:59 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.794	152	2778	0.400	ng/ul	0.00
4) Naphthalene-d8	10.581	136	11969	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.421	164	7016	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.159	188	12749	0.400	ng/ul	0.00
17) Chrysene-d12	21.323	240	11339	0.400	ng/ul	0.00
23) Perylene-d12	23.576	264	10239	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.283	96	1107	0.372	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.169	152	8385	0.438	ng/ul	0.00
18) Fluoranthene-d10	19.180	212	12551	0.361	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.317	88	1143	0.369	ng/ul	97
5) Naphthalene	10.631	128	14822	0.416	ng/ul	99
7) 2-Methylnaphthalene	12.245	142	10613	0.442	ng/ul	99
8) 1-Methylnaphthalene	12.461	142	9856	0.422	ng/ul	99
10) Acenaphthylene	14.140	152	13945	0.425	ng/ul	99
11) Acenaphthene	14.482	153	11016	0.412	ng/ul	99
12) Fluorene	15.471	166	12155	0.410	ng/ul	99
14) Pentachlorophenol	16.826	266	1259	0.339	ng/ul	99
15) Phenanthrene	17.200	178	18290	0.403	ng/ul	100
16) Anthracene	17.294	178	17530	0.441	ng/ul	100
19) Fluoranthene	19.211	202	21779	0.418	ng/ul	98
20) Pyrene	19.573	202	21572	0.404	ng/ul	98
21) Benzo(a)anthracene	21.307	228	17523	0.389	ng/ul	99
22) Chrysene	21.357	228	20355	0.407	ng/ul	99
24) Benzo(b)fluoranthene	22.895	252	18595	0.417	ng/ul	98
25) Benzo(k)fluoranthene	22.941	252	18425	0.394	ng/ul	98
26) Benzo(a)pyrene	23.477	252	15978	0.404	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.856	276	20829	0.414	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.868	278	16978	0.424	ng/ul	99
29) Benzo(g,h,i)perylene	26.553	276	17943	0.413	ng/ul	97

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

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