

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110921\
 Data File : BM032935.D
 Acq On : 10 Nov 2021 04:00
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4162

Quant Time: Nov 10 06:11:35 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM110921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 09 13:31:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.466	152	2110	0.400	ng/ul	0.00
4) Naphthalene-d8	10.239	136	7541	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.121	164	4428	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.865	188	9259	0.400	ng/ul	0.00
17) Chrysene-d12	21.050	240	7926	0.400	ng/ul	0.00
23) Perylene-d12	23.159	264	6545	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.867	96	1082	0.402	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.840	152	4190	0.391	ng/ul	0.00
18) Fluoranthene-d10	18.892	212	9722	0.405	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.902	88	1191	0.430	ng/ul	95
5) Naphthalene	10.289	128	9252	0.423	ng/ul	99
7) 2-Methylnaphthalene	11.912	142	5954	0.393	ng/ul	98
8) 1-Methylnaphthalene	12.137	142	5878	0.396	ng/ul	100
10) Acenaphthylene	13.842	152	7512	0.390	ng/ul	100
11) Acenaphthene	14.186	153	6838	0.407	ng/ul	99
12) Fluorene	15.175	166	7790	0.393	ng/ul	99
14) Pentachlorophenol	16.539	266	829	0.344	ng/ul	99
15) Phenanthrene	16.904	178	12447	0.405	ng/ul	100
16) Anthracene	16.997	178	10571	0.391	ng/ul	100
19) Fluoranthene	18.923	202	14644	0.394	ng/ul	97
20) Pyrene	19.285	202	15034	0.389	ng/ul	98
21) Benzo(a)anthracene	21.035	228	11416	0.390	ng/ul	100
22) Chrysene	21.086	228	13326	0.404	ng/ul	99
24) Benzo(b)fluoranthene	22.534	252	12170	0.405	ng/ul	99
25) Benzo(k)fluoranthene	22.575	252	13282	0.413	ng/ul	99
26) Benzo(a)pyrene	23.067	252	9965	0.403	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.218	276	11784	0.414	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.221	278	9362	0.415	ng/ul	99
29) Benzo(g,h,i)perylene	25.851	276	10392	0.420	ng/ul	100

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

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