

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102122\
 Data File : BM037180.D
 Acq On : 21 Oct 2022 22:44
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4158

Quant Time: Oct 22 03:35:02 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM101922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 22 03:34:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.900	152	4770	0.400	ng/ul	0.00
4) Naphthalene-d8	10.699	136	14392	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.530	164	8315	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.272	188	17951	0.400	ng/ul	0.00
17) Chrysene-d12	21.446	240	14084	0.400	ng/ul	0.00
23) Perylene-d12	23.823	264	11546	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.314	96	2318	0.399	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.288	152	8433	0.412	ng/ul	0.00
18) Fluoranthene-d10	19.295	212	18410	0.432	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	2221	0.368	ng/ul#	92
5) Naphthalene	10.748	128	16869	0.396	ng/ul	99
7) 2-Methylnaphthalene	12.360	142	10561	0.414	ng/ul	97
8) 1-Methylnaphthalene	12.580	142	10854	0.417	ng/ul	99
10) Acenaphthylene	14.252	152	14358	0.402	ng/ul	99
11) Acenaphthene	14.594	153	12521	0.388	ng/ul	98
12) Fluorene	15.580	166	14299	0.379	ng/ul	98
14) Pentachlorophenol	16.926	266	3049	0.702	ng/ul	99
15) Phenanthrene	17.314	178	23655	0.390	ng/ul	99
16) Anthracene	17.403	178	19352	0.403	ng/ul	99
19) Fluoranthene	19.327	202	26468	0.433	ng/ul	98
20) Pyrene	19.685	202	26929	0.434	ng/ul	97
21) Benzo(a)anthracene	21.432	228	22410	0.410	ng/ul	99
22) Chrysene	21.484	228	24618	0.391	ng/ul	100
24) Benzo(b)fluoranthene	23.095	252	24479	0.400	ng/ul	99
25) Benzo(k)fluoranthene	23.144	252	22727	0.385	ng/ul	100
26) Benzo(a)pyrene	23.712	252	19468	0.396	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.275	276	22759	0.332	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.292	278	17901	0.331	ng/ul	100
29) Benzo(g,h,i)perylene	27.026	276	19239	0.319	ng/ul	100

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

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