

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100722\
 Data File : BM036866.D
 Acq On : 07 Oct 2022 20:41
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4134

Quant Time: Oct 08 02:50:30 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM100622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 06 17:02:30 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.968	152	5929	0.400	ng/ul	0.00
4) Naphthalene-d8	10.770	136	20185	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.590	164	10918	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.327	188	22474	0.400	ng/ul	0.00
17) Chrysene-d12	21.496	240	19440	0.400	ng/ul	0.00
23) Perylene-d12	23.899	264	16560	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.356	96	3192	0.382	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.360	152	11983	0.393	ng/ul	0.00
18) Fluoranthene-d10	19.346	212	23176	0.399	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.390	88	3218	0.394	ng/ul#	85
5) Naphthalene	10.825	128	22915	0.397	ng/ul	97
7) 2-Methylnaphthalene	12.431	142	14128	0.400	ng/ul	92
8) 1-Methylnaphthalene	12.646	142	14341	0.400	ng/ul	99
10) Acenaphthylene	14.312	152	17756	0.391	ng/ul	99
11) Acenaphthene	14.654	153	15798	0.401	ng/ul	99
12) Fluorene	15.635	166	17718	0.397	ng/ul	99
14) Pentachlorophenol	16.981	266	2040	0.375	ng/ul	97
15) Phenanthrene	17.369	178	28703	0.400	ng/ul	99
16) Anthracene	17.462	178	23609	0.394	ng/ul	98
19) Fluoranthene	19.378	202	30796	0.398	ng/ul	99
20) Pyrene	19.736	202	31657	0.393	ng/ul	98
21) Benzo(a)anthracene	21.478	228	27369	0.390	ng/ul	99
22) Chrysene	21.531	228	29992	0.398	ng/ul	98
24) Benzo(b)fluoranthene	23.162	252	30453	0.412	ng/ul	94
25) Benzo(k)fluoranthene	23.212	252	28633	0.397	ng/ul	92
26) Benzo(a)pyrene	23.790	252	24932	0.415	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.392	276	33448	0.393	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.409	278	26575	0.387	ng/ul#	90
29) Benzo(g,h,i)perylene	27.160	276	28022	0.374	ng/ul	98

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

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