

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102522\
 Data File : BM037220.D
 Acq On : 25 Oct 2022 10:02
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4163

Quant Time: Oct 26 07:31:24 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	5879	0.400	ng/u1	0.00
4) Naphthalene-d8	10.699	136	18019	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.529	164	9883	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.271	188	18252	0.400	ng/u1	0.00
17) Chrysene-d12	21.447	240	10940	0.400	ng/u1	0.00
23) Perylene-d12	23.818	264	10962	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.314	96	2635	0.368	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.283	152	9897	0.386	ng/u1	0.00
18) Fluoranthene-d10	19.294	212	14737	0.445	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.352	88	2575	0.346	ng/u1	94
5) Naphthalene	10.748	128	19715	0.369	ng/u1	99
7) 2-Methylnaphthalene	12.360	142	12379	0.387	ng/u1	97
8) 1-Methylnaphthalene	12.580	142	12605	0.386	ng/u1	99
10) Acenaphthylene	14.251	152	16456	0.388	ng/u1	100
11) Acenaphthene	14.589	153	13698	0.357	ng/u1	98
12) Fluorene	15.574	166	14956	0.333	ng/u1	99
14) Pentachlorophenol	16.925	266	1489	0.337	ng/u1	99
15) Phenanthrene	17.309	178	22142	0.359	ng/u1	99
16) Anthracene	17.402	178	18312	0.375	ng/u1	99
19) Fluoranthene	19.322	202	20799	0.438	ng/u1	100
20) Pyrene	19.689	202	20792	0.432	ng/u1	98
21) Benzo(a)anthracene	21.430	228	15925	0.375	ng/u1	100
22) Chrysene	21.482	228	17314	0.354	ng/u1	99
24) Benzo(b)fluoranthene	23.093	252	19064	0.328	ng/u1	98
25) Benzo(k)fluoranthene	23.139	252	18282	0.326	ng/u1	98
26) Benzo(a)pyrene	23.709	252	16734	0.359	ng/u1	99
27) Indeno(1,2,3-cd)pyrene	26.269	276	22810	0.350	ng/u1#	96
28) Dibenzo(a,h)anthracene	26.286	278	17792	0.346	ng/u1	98
29) Benzo(g,h,i)perylene	27.027	276	19405	0.339	ng/u1	98

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

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