

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102122\
 Data File : BM037171.D
 Acq On : 21 Oct 2022 13:24
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4157

Quant Time: Oct 22 02:32:10 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 : Thu Oct 20 02:11:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.909	152	4193	0.400	ng/u1	0.00
4) Naphthalene-d8	10.705	136	12782	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.533	164	7502	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.275	188	16603	0.400	ng/u1	0.00
17) Chrysene-d12	21.453	240	13614	0.400	ng/u1	0.00
23) Perylene-d12	23.830	264	10530	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	1961	0.384	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.294	152	7501	0.413	ng/u1	0.00
18) Fluoranthene-d10	19.299	212	17256	0.419	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.352	88	2000	0.377	ng/u1	92
5) Naphthalene	10.754	128	14933	0.394	ng/u1	100
7) 2-Methylnaphthalene	12.365	142	9391	0.414	ng/u1	98
8) 1-Methylnaphthalene	12.585	142	9589	0.414	ng/u1	100
10) Acenaphthylene	14.256	152	13158	0.408	ng/u1	99
11) Acenaphthene	14.598	153	11318	0.389	ng/u1	98
12) Fluorene	15.584	166	13100	0.385	ng/u1	99
14) Pentachlorophenol	16.933	266	1631	0.406	ng/u1	98
15) Phenanthrene	17.318	178	21608	0.386	ng/u1	99
16) Anthracene	17.410	178	17974	0.405	ng/u1	99
19) Fluoranthene	19.331	202	24439	0.413	ng/u1	98
20) Pyrene	19.694	202	25053	0.418	ng/u1	100
21) Benzo(a)anthracene	21.436	228	21370	0.404	ng/u1	99
22) Chrysene	21.488	228	23376	0.385	ng/u1	99
24) Benzo(b)fluoranthene	23.104	252	21951	0.394	ng/u1	99
25) Benzo(k)fluoranthene	23.151	252	21018	0.390	ng/u1	99
26) Benzo(a)pyrene	23.724	252	16846	0.376	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	26.289	276	19665	0.314	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.306	278	15671	0.317	ng/u1	98
29) Benzo(g,h,i)perylene	27.041	276	16963	0.308	ng/u1	100

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

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