

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101521\
 Data File : BM032515.D
 Acq On : 15 Oct 2021 08:38
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4137

Quant Time: Oct 15 12:35:52 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM101421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 14 19:35:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.556	152	4263	0.40	ng/ul	0.00
4) Naphthalene-d8	10.344	136	12957	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.217	164	6809	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.953	188	13584	0.40	ng/ul	0.00
17) Chrysene-d12	21.135	240	9323	0.40	ng/ul	0.00
23) Perylene-d12	23.278	264	7062	0.40	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.927	96	2024	0.42	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.940	152	6743	0.39	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	12799	0.44	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.961	88	2198	0.42	ng/ul#	86
5) Naphthalene	10.393	128	15530	0.40	ng/ul	99
7) 2-Methylnaphthalene	12.012	142	10069	0.39	ng/ul	100
8) 1-Methylnaphthalene	12.237	142	10032	0.40	ng/ul	99
10) Acenaphthylene	13.933	152	10982	0.38	ng/ul	100
11) Acenaphthene	14.278	153	10894	0.41	ng/ul	99
12) Fluorene	15.263	166	11970	0.40	ng/ul	99
14) Pentachlorophenol	16.623	266	1239	0.35	ng/ul	98
15) Phenanthrene	16.995	178	18682	0.41	ng/ul	99
16) Anthracene	17.085	178	15361	0.38	ng/ul	99
19) Fluoranthene	19.011	202	20358	0.44	ng/ul	99
20) Pyrene	19.374	202	20206	0.43	ng/ul	100
21) Benzo(a)anthracene	21.121	228	14117	0.39	ng/ul	99
22) Chrysene	21.171	228	16517	0.41	ng/ul	100
24) Benzo(b)fluoranthene	22.643	252	14604	0.45	ng/ul	95
25) Benzo(k)fluoranthene	22.685	252	14382	0.44	ng/ul	97
26) Benzo(a)pyrene	23.184	252	11370	0.42	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	25.386	276	13599	0.41	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.393	278	10638	0.40	ng/ul	98
29) Benzo(g,h,i)perylene	26.031	276	12173	0.42	ng/ul	98

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

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