

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080221\
 Data File : BM031440.D
 Acq On : 04 Aug 2021 09:11
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4092

Quant Time: Aug 04 12:16:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 04 11:34:03 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.597	152	1035	0.400	ng/ul	0.00
4) Naphthalene-d8	10.355	136	3758	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.227	164	2160	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.974	188	4186	0.400	ng/ul	0.00
17) Chrysene-d12	21.168	240	3267	0.400	ng/ul	0.00
23) Perylene-d12	23.331	264	3139	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	416	0.362	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.952	152	2194	0.346	ng/ul	0.00
18) Fluoranthene-d10	19.008	212	3881	0.356	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.231	88	404	0.348	ng/ul	90
5) Naphthalene	10.405	128	3992	0.357	ng/ul	98
7) 2-Methylnaphthalene	12.028	142	2704	0.354	ng/ul	98
8) 1-Methylnaphthalene	12.249	142	2603	0.344	ng/ul	100
10) Acenaphthylene	13.949	152	3327	0.362	ng/ul	100
11) Acenaphthene	14.288	153	2692	0.348	ng/ul	98
12) Fluorene	15.281	166	2983	0.335	ng/ul	99
14) Pentachlorophenol	16.631	266	437	0.326	ng/ul	96
15) Phenanthrene	17.016	178	4548	0.342	ng/ul	100
16) Anthracene	17.110	178	4245	0.341	ng/ul	98
19) Fluoranthene	19.039	202	5015	0.359	ng/ul	99
20) Pyrene	19.401	202	5117	0.362	ng/ul	98
21) Benzo(a)anthracene	21.154	228	4531	0.343	ng/ul	99
22) Chrysene	21.204	228	4771	0.350	ng/ul	99
24) Benzo(b)fluoranthene	22.686	252	4821	0.341	ng/ul	96
25) Benzo(k)fluoranthene	22.727	252	4869	0.333	ng/ul	96
26) Benzo(a)pyrene	23.234	252	4222	0.341	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	25.465	276	5381	0.336	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.485	278	4256	0.330	ng/ul	97
29) Benzo(g,h,i)perylene	26.119	276	4642	0.340	ng/ul	96

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

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