

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012022\
 Data File : BM033815.D
 Acq On : 21 Jan 2022 08:00
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4197

Quant Time: Jan 21 08:37:32 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM010422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 21 01:34:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.949	152	3219	0.400	ng/ul	0.00
4) Naphthalene-d8	10.760	136	11113	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.583	164	5426	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.319	188	9787	0.400	ng/ul	0.00
17) Chrysene-d12	21.482	240	7118	0.400	ng/ul #	0.00
23) Perylene-d12	23.880	264	7071	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.303	96	1413	0.412	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.348	152	5995	0.397	ng/ul	0.00
18) Fluoranthene-d10	19.337	212	9168	0.416	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.337	88	1406	0.376	ng/ul#	84
5) Naphthalene	10.810	128	11973	0.372	ng/ul#	88
7) 2-Methylnaphthalene	12.420	142	7772	0.376	ng/ul	99
8) 1-Methylnaphthalene	12.636	142	7632	0.376	ng/ul	99
10) Acenaphthylene	14.306	152	9440	0.376	ng/ul#	90
11) Acenaphthene	14.648	153	7627	0.381	ng/ul	95
12) Fluorene	15.630	166	8272	0.346	ng/ul#	95
14) Pentachlorophenol	16.976	266	1394	0.361	ng/ul	98
15) Phenanthrene	17.361	178	12150	0.383	ng/ul	95
16) Anthracene	17.455	178	11083	0.381	ng/ul	93
19) Fluoranthene	19.368	202	13542	0.406	ng/ul	100
20) Pyrene	19.726	202	13767	0.408	ng/ul#	94
21) Benzo(a)anthracene	21.465	228	11243	0.380	ng/ul	98
22) Chrysene	21.518	228	12057	0.399	ng/ul	98
24) Benzo(b)fluoranthene	23.148	252	12513	0.404	ng/ul	91
25) Benzo(k)fluoranthene	23.194	252	12325	0.406	ng/ul#	91
26) Benzo(a)pyrene	23.773	252	11029	0.410	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	26.387	276	14707	0.419	ng/ul#	87
28) Dibenzo(a,h)anthracene	26.404	278	11711	0.420	ng/ul	92
29) Benzo(g,h,i)perylene	27.151	276	12579	0.417	ng/ul#	93

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

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