

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012422\
 Data File : BM033838.D
 Acq On : 24 Jan 2022 15:27
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV049

Quant Time: Jan 24 16:55:39 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM012422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 24 15:20:03 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.943	152	2072	0.400	ng/ul	0.00
4) Naphthalene-d8	10.756	136	7120	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.583	164	3570	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.319	188	7318	0.400	ng/ul	0.00
17) Chrysene-d12	21.485	240	5977	0.400	ng/ul #	0.00
23) Perylene-d12	23.887	264	5863	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.296	96	965	0.446	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.344	152	4052	0.413	ng/ul	0.00
18) Fluoranthene-d10	19.342	212	7616	0.414	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.334	88	976	0.430	ng/ul#	88
5) Naphthalene	10.805	128	7965	0.407	ng/ul#	89
7) 2-Methylnaphthalene	12.416	142	5238	0.412	ng/ul	100
8) 1-Methylnaphthalene	12.636	142	5092	0.408	ng/ul	100
10) Acenaphthylene	14.303	152	6329	0.397	ng/ul#	92
11) Acenaphthene	14.644	153	5301	0.415	ng/ul	96
12) Fluorene	15.626	166	5839	0.401	ng/ul	95
14) Pentachlorophenol	16.972	266	1401	0.417	ng/ul	96
15) Phenanthrene	17.361	178	9470	0.406	ng/ul	95
16) Anthracene	17.455	178	8670	0.397	ng/ul	92
19) Fluoranthene	19.368	202	11229	0.412	ng/ul	99
20) Pyrene	19.731	202	11461	0.411	ng/ul	95
21) Benzo(a)anthracene	21.470	228	9936	0.407	ng/ul	97
22) Chrysene	21.521	228	10531	0.407	ng/ul	97
24) Benzo(b)fluoranthene	23.153	252	10811	0.411	ng/ul	92
25) Benzo(k)fluoranthene	23.202	252	10813	0.410	ng/ul#	92
26) Benzo(a)pyrene	23.781	252	9225	0.404	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.395	276	12434	0.409	ng/ul#	87
28) Dibenzo(a,h)anthracene	26.427	278	9778	0.408	ng/ul	92
29) Benzo(g,h,i)perylene	27.163	276	10658	0.413	ng/ul	93

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

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