

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062821\
 Data File : BM030655.D
 Acq On : 28 Jun 2021 14:38
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV710

Quant Time: Jun 28 15:23:01 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 28 14:33:12 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.791	152	2064	0.400	ng/ul	0.00
4) Naphthalene-d8	10.576	136	7218	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.421	164	3798	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.158	188	6994	0.400	ng/ul	0.00
17) Chrysene-d12	21.328	240	6704	0.400	ng/ul	0.00
23) Perylene-d12	23.585	264	6285	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.282	96	855	0.380	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.169	152	4308	0.380	ng/ul	0.00
18) Fluoranthene-d10	19.180	212	7437	0.387	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.317	88	861	0.371	ng/ul	88
5) Naphthalene	10.626	128	7519	0.366	ng/ul	100
7) 2-Methylnaphthalene	12.241	142	5014	0.362	ng/ul	99
8) 1-Methylnaphthalene	12.461	142	4589	0.341	ng/ul	99
10) Acenaphthylene	14.140	152	6366	0.375	ng/ul	99
11) Acenaphthene	14.481	153	5010	0.365	ng/ul	99
12) Fluorene	15.467	166	5448	0.359	ng/ul	99
14) Pentachlorophenol	16.822	266	739	0.324	ng/ul	97
15) Phenanthrene	17.200	178	8574	0.365	ng/ul	99
16) Anthracene	17.290	178	7186	0.349	ng/ul	99
19) Fluoranthene	19.210	202	10525	0.356	ng/ul	99
20) Pyrene	19.573	202	10525	0.354	ng/ul	99
21) Benzo(a)anthracene	21.314	228	9046	0.338	ng/ul	99
22) Chrysene	21.364	228	10788	0.360	ng/ul	99
24) Benzo(b)fluoranthene	22.907	252	10688	0.384	ng/ul	99
25) Benzo(k)fluoranthene	22.951	252	10613	0.370	ng/ul	99
26) Benzo(a)pyrene	23.486	252	9496	0.401	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.868	276	11675	0.379	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.880	278	9424	0.381	ng/ul	100
29) Benzo(g,h,i)perylene	26.565	276	10461	0.390	ng/ul	98

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

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