

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012422\
 Data File : BM033836.D
 Acq On : 24 Jan 2022 13:54
 Operator : CG/JU
 Sample : SSTD1.647
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.6047

Quant Time: Jan 24 15:12:17 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 14:03:45 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	2116	0.400	ng/ul	0.00
4) Naphthalene-d8	10.756	136	7286	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.579	164	3839	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	7723	0.400	ng/ul	0.00
17) Chrysene-d12	21.480	240	6339	0.400	ng/ul #	0.00
23) Perylene-d12	23.883	264	6161	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.296	96	3393	1.505	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.344	152	15595	1.574	ng/ul	0.00
18) Fluoranthene-d10	19.338	212	29656	1.510	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.334	88	3530	1.436	ng/ul#	80
5) Naphthalene	10.805	128	30215	1.431	ng/ul#	87
7) 2-Methylnaphthalene	12.415	142	20170	1.488	ng/ul	100
8) 1-Methylnaphthalene	12.636	142	19722	1.481	ng/ul	99
10) Acenaphthylene	14.303	152	25480	1.436	ng/ul#	89
11) Acenaphthene	14.644	153	20536	1.451	ng/ul	94
12) Fluorene	15.626	166	23482	1.387	ng/ul#	96
14) Pentachlorophenol	16.972	266	5270	1.729	ng/ul	98
15) Phenanthrene	17.358	178	36937	1.477	ng/ul	93
16) Anthracene	17.448	178	34787	1.516	ng/ul#	91
19) Fluoranthene	19.368	202	43833	1.474	ng/ul	98
20) Pyrene	19.727	202	43915	1.460	ng/ul	95
21) Benzo(a)anthracene	21.465	228	38898	1.477	ng/ul	97
22) Chrysene	21.518	228	40308	1.497	ng/ul	96
24) Benzo(b)fluoranthene	23.148	252	41348	1.532	ng/ul	88
25) Benzo(k)fluoranthene	23.194	252	42458	1.604	ng/ul#	86
26) Benzo(a)pyrene	23.776	252	36333	1.548	ng/ul#	80
27) Indeno(1,2,3-cd)pyrene	26.385	276	47807	1.565	ng/ul#	85
28) Dibenzo(a,h)anthracene	26.410	278	37933	1.562	ng/ul#	85
29) Benzo(g,h,i)perylene	27.158	276	40648	1.548	ng/ul	94

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

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