

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030422\
 Data File : BM034133.D
 Acq On : 04 Mar 2022 12:20
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4059

Quant Time: Mar 04 13:48:08 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM030322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 03 14:59:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.001	152	2598	0.400	ng/ul	0.00
4) Naphthalene-d8	10.812	136	6744	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.634	164	4313	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.371	188	9624	0.400	ng/ul #	0.00
17) Chrysene-d12	21.542	240	7981	0.400	ng/ul #	0.00
23) Perylene-d12	23.988	264	7229	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.377	96	1263	0.418	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.396	152	4188	0.403	ng/ul	0.00
18) Fluoranthene-d10	19.394	212	9966	0.414	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.410	88	1207	0.412	ng/ul#	71
5) Naphthalene	10.862	128	7634	0.412	ng/ul#	89
7) 2-Methylnaphthalene	12.467	142	5386	0.412	ng/ul	99
8) 1-Methylnaphthalene	12.687	142	5309	0.408	ng/ul	98
10) Acenaphthylene	14.352	152	7490	0.412	ng/ul#	90
11) Acenaphthene	14.694	153	6094	0.414	ng/ul	95
12) Fluorene	15.675	166	7275	0.408	ng/ul#	95
14) Pentachlorophenol	17.021	266	1141	0.370	ng/ul	97
15) Phenanthrene	17.413	178	11873	0.409	ng/ul	94
16) Anthracene	17.506	178	10656	0.404	ng/ul	92
19) Fluoranthene	19.422	202	13587	0.415	ng/ul#	89
20) Pyrene	19.785	202	13666	0.417	ng/ul#	90
21) Benzo(a)anthracene	21.527	228	11421	0.411	ng/ul	96
22) Chrysene	21.580	228	12132	0.411	ng/ul	98
24) Benzo(b)fluoranthene	23.240	252	11928	0.405	ng/ul	90
25) Benzo(k)fluoranthene	23.290	252	12356	0.413	ng/ul#	88
26) Benzo(a)pyrene	23.877	252	10236	0.412	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	26.536	276	14075	0.405	ng/ul#	57
28) Dibenzo(a,h)anthracene	26.562	278	11323	0.399	ng/ul#	86
29) Benzo(g,h,i)perylene	27.317	276	12444	0.408	ng/ul#	91

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

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