

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040623\
 Data File : BM039315.D
 Acq On : 06 Apr 2023 13:33
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV021

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 04/07/2023
 Supervised By :Jagrut Upadhyay 04/07/2023

Quant Time: Apr 06 23:59:41 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM040623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 06 23:58:10 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	3580	0.400	ng/ul	0.00
4) Naphthalene-d8	10.702	136	17468	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.537	164	10231	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.291	188	15621	0.400	ng/ul	0.00
17) Chrysene-d12	21.480	240	13779	0.400	ng/ul	0.00
23) Perylene-d12	23.868	264	14689	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.263	96	3396	0.512	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.297	152	9503	0.471	ng/ul	0.00
18) Fluoranthene-d10	19.320	212	19507	0.454	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.296	88	3643	0.513	ng/ul#	90
5) Naphthalene	10.752	128	19032	0.458	ng/ul	99
7) 2-Methylnaphthalene	12.368	142	10673	0.457	ng/ul	100
8) 1-Methylnaphthalene	12.583	142	12177	0.461	ng/ul	98
10) Acenaphthylene	14.259	152	18164	0.458	ng/ul	99
11) Acenaphthene	14.601	153	14650	0.469	ng/ul	99
12) Fluorene	15.591	166	14589	0.469	ng/ul	99
14) Pentachlorophenol	16.957	266	1552	0.417	ng/ul#	81
15) Phenanthrene	17.329	178	19887	0.459	ng/ul	99
16) Anthracene	17.426	178	14438	0.438	ng/ul	99
19) Fluoranthene	19.348	202	25577	0.448	ng/ul	100
20) Pyrene	19.710	202	29943	0.458	ng/ul	100
21) Benzo(a)anthracene	21.463	228	13603m	0.428	ng/ul	
22) Chrysene	21.513	228	19183	0.492	ng/ul	99
24) Benzo(b)fluoranthene	23.140	252	21139	0.454	ng/ul	94
25) Benzo(k)fluoranthene	23.187	252	19301	0.434	ng/ul	93
26) Benzo(a)pyrene	23.766	252	20771	0.442	ng/ul	91
27) Indeno(1,2,3-cd)pyrene	26.357	276	24276	0.410	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.377	278	16585	0.401	ng/ul	98
29) Benzo(g,h,i)perylene	27.109	276	23447	0.413	ng/ul	98

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

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