

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050823\
 Data File : BN025339.D
 Acq On : 08 May 2023 18:57
 Operator : CG/JU
 Sample : 02588-08MS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ZONE-A-3-6-05022023MS

Quant Time: May 09 04:03:35 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN050623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 08 18:27:02 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 05/09/2023
 Supervised By :Jagrut Upadhyay 05/09/2023

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.903	152	5292	0.400	ng	0.00
7) Naphthalene-d8	10.707	136	17880	0.400	ng	#-0.01
13) Acenaphthene-d10	14.544	164	10738	0.400	ng	0.00
19) Phenanthrene-d10	17.298	188	24721	0.400	ng	0.00
29) Chrysene-d12	21.494	240	26135	0.400	ng	0.00
35) Perylene-d12	23.914	264	26794	0.400	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.462	112	3841	0.298	ng	0.00
5) Phenol-d6	7.080	99	5162	0.319	ng	0.00
8) Nitrobenzene-d5	9.073	82	3842	0.261	ng	-0.01
11) 2-Methylnaphthalene-d10	12.301	152	10465m	0.428	ng	0.00
14) 2,4,6-Tribromophenol	16.044	330	1487	0.271	ng	-0.01
15) 2-Fluorobiphenyl	13.175	172	10038	0.273	ng	0.00
27) Fluoranthene-d10	19.329	212	16018	0.262	ng	0.00
31) Terphenyl-d14	19.923	244	12218	0.232	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.310	88	1682	0.303	ng	# 56
3) n-Nitrosodimethylamine	3.642	42	3191	0.376	ng	# 98
6) bis(2-Chloroethyl)ether	7.326	93	6136	0.431	ng	96
9) Naphthalene	10.760	128	32825	0.716	ng	99
10) Hexachlorobutadiene	11.038	225	3295	0.380	ng	# 97
12) 2-Methylnaphthalene	12.373	142	17959	0.554	ng	99
16) Acenaphthylene	14.266	152	39859	0.829	ng	99
17) Acenaphthene	14.608	154	16577	0.554	ng	99
18) Fluorene	15.602	166	22688	0.554	ng	98
20) 4,6-Dinitro-2-methylph...	15.709	198	322	0.340	ng	# 75
21) 4-Bromophenyl-phenylether	16.491	248	5146	0.373	ng	# 78
22) Hexachlorobenzene	16.603	284	5966	0.401	ng	99
23) Atrazine	16.752	200	4731	0.392	ng	100
24) Pentachlorophenol	16.963	266	2906	0.771	ng	91
25) Phenanthrene	17.335	178	152775	2.234	ng	100
26) Anthracene	17.435	178	55495	0.866	ng	99
28) Fluoranthene	19.359	202	364859	4.269	ng	99
30) Pyrene	19.725	202	341424	3.821	ng	97
32) Benzo(a)anthracene	21.477	228	212440	2.405	ng	99
33) Chrysene	21.530	228	227181	2.584	ng	98
34) Bis(2-ethylhexyl)phtha...	21.387	149	43235	0.624	ng	99
36) Indeno(1,2,3-cd)pyrene	26.431	276	136858	1.274	ng	96
37) Benzo(b)fluoranthene	23.177	252	281460	3.314	ng	95
38) Benzo(k)fluoranthene	23.221	252	129494m	1.456	ng	
39) Benzo(a)pyrene	23.803	252	202143	2.323	ng	93
40) Dibenzo(a,h)anthracene	26.440	278	52323	0.617	ng	97
41) Benzo(g,h,i)perylene	27.206	276	126735	1.360	ng	99

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

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