

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050623\
 Data File : BN025319.D
 Acq On : 06 May 2023 17:08
 Operator : CG/JU
 Sample : 02588-10
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ZONE-D-0-3-05022023

Quant Time: May 08 04:01:21 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 02:37:05 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 05/08/2023
 Supervised By : Jagrut Upadhyay 05/08/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.911	152	9996	0.400	ng	0.00
7) Naphthalene-d8	10.718	136	31673	0.400	ng	-0.01
13) Acenaphthene-d10	14.544	164	19110	0.400	ng	-0.01
19) Phenanthrene-d10	17.298	188	39805	0.400	ng	#-0.01
29) Chrysene-d12	21.495	240	27038	0.400	ng	#-0.01
35) Perylene-d12	23.940	264	26866	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.470	112	6557	0.269	ng	0.00
5) Phenol-d6	7.080	99	8755	0.287	ng	0.00
8) Nitrobenzene-d5	9.073	82	6342	0.243	ng	-0.02
11) 2-Methylnaphthalene-d10	12.305	152	11946	0.276	ng	-0.01
14) 2,4,6-Tribromophenol	16.045	330	2341	0.240	ng	-0.02
15) 2-Fluorobiphenyl	13.176	172	17626	0.269	ng	-0.01
27) Fluoranthene-d10	19.334	212	21182	0.215	ng	0.00
31) Terphenyl-d14	19.924	244	18138	0.333	ng	-0.01
Target Compounds						
9) Naphthalene	10.760	128	58398	0.719	ng	100
12) 2-Methylnaphthalene	12.377	142	25633	0.446	ng	99
16) Acenaphthylene	14.266	152	76952	0.899	ng	99
17) Acenaphthene	14.608	154	79111	1.487	ng	100
18) Fluorene	15.603	166	102071	1.400	ng	# 98
25) Phenanthrene	17.336	178	1640911	14.900	ng	100
26) Anthracene	17.435	178	423395	4.104	ng	# 97
28) Fluoranthene	19.367	202	3065085	22.274	ng	99
30) Pyrene	19.730	202	2580369	27.916	ng	# 94
32) Benzo(a)anthracene	21.477	228	1215484	13.300	ng	98
33) Chrysene	21.531	228	1124118	12.357	ng	97
34) Bis(2-ethylhexyl)phtha...	21.396	149	90237	1.258	ng	93
36) Indeno(1,2,3-cd)pyrene	26.487	276	623769	5.792	ng	96
37) Benzo(b)fluoranthene	23.198	252	1331120	15.630	ng	95
38) Benzo(k)fluoranthene	23.239	252	493137m	5.532	ng	
39) Benzo(a)pyrene	23.835	252	955602m	10.953	ng	
40) Dibenzo(a,h)anthracene	26.496	278	158286	1.861	ng	# 89
41) Benzo(g,h,i)perylene	27.273	276	598168	6.404	ng	100

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

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