

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032122\
 Data File : BN019027.D
 Acq On : 21 Mar 2022 17:54
 Operator : CG/JU
 Sample : N2007-01 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SP1

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/22/2022
 Supervised By : Jagrut Upadhyay 03/23/2022

Quant Time: Mar 22 02:31:20 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 17 13:18:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.854	152	7181	0.400	ng	0.00
7) Naphthalene-d8	10.658	136	19658	0.400	ng	# 0.00
13) Acenaphthene-d10	14.495	164	11564	0.400	ng	0.00
19) Phenanthrene-d10	17.226	188	21094	0.400	ng	-0.01
29) Chrysene-d12	21.413	240	13743	0.400	ng	# 0.00
35) Perylene-d12	23.782	264	12933	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.406	112	776	0.048	ng	0.00
5) Phenol-d6	7.024	99	595	0.035	ng	0.01
8) Nitrobenzene-d5	9.025	82	447	0.032	ng	0.01
11) 2-Methylnaphthalene-d10	12.253	152	1357	0.042	ng	0.00
14) 2,4,6-Tribromophenol	15.985	330	204	0.037	ng	0.00
15) 2-Fluorobiphenyl	13.127	172	2042	0.044	ng	0.00
27) Fluoranthene-d10	19.261	212	2343	0.037	ng	0.00
31) Terphenyl-d14	19.860	244	1754	0.052	ng	0.00
Target Compounds						Qvalue
9) Naphthalene	10.712	128	3126	0.056	ng	# 94
12) 2-Methylnaphthalene	12.325	142	2421	0.064	ng	94
16) Acenaphthylene	14.207	152	5371	0.100	ng	99
17) Acenaphthene	14.549	154	9001	0.239	ng	99
18) Fluorene	15.532	166	11180	0.236	ng	98
25) Phenanthrene	17.267	178	191762	2.907	ng	100
26) Anthracene	17.359	178	30095	0.540	ng	97
28) Fluoranthene	19.291	202	270276	3.492	ng	99
30) Pyrene	19.653	202	228842	3.343	ng	97
32) Benzo(a)anthracene	21.396	228	79739	1.883	ng	98
33) Chrysene	21.449	228	84879	1.486	ng	98
34) Bis(2-ethylhexyl)phtha...	21.333	149	9832	0.350	ng	# 99
36) Indeno(1,2,3-cd)pyrene	26.232	276	37079	0.795	ng	# 88
37) Benzo(b)fluoranthene	23.065	252	90676	1.922	ng	96
38) Benzo(k)fluoranthene	23.106	252	33955m	0.525	ng	
39) Benzo(a)pyrene	23.679	252	57163m	1.332	ng	
40) Dibenzo(a,h)anthracene	26.241	278	7472	0.206	ng	# 79
41) Benzo(g,h,i)perylene	26.983	276	44796	0.978	ng	99

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032122\
Data File : BN019027.D
Acq On : 21 Mar 2022 17:54
Operator : CG/JU
Sample : N2007-01 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SP1

Quant Time: Mar 22 02:31:20 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN031722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 17 13:18:23 2022
Response via : Initial Calibration

Manual Integrations
APPROVED

Reviewed By :Christian Giraldo 03/22/2022
Supervised By :Jagrut Upadhyay 03/23/2022

