

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN113021\
 Data File : BN017646.D
 Acq On : 01 Dec 2021 12:10
 Operator : CG/JU
 Sample : M4862-08
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CS-4A

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/01/2021
 Supervised By :mohammad ahmed 12/05/2021

Quant Time: Dec 01 12:42:21 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN113021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 30 18:26:44 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.773	152	21130	0.400	ng	0.00
7) Naphthalene-d8	10.548	136	89779	0.400	ng	# 0.00
13) Acenaphthene-d10	14.384	164	57087	0.400	ng	0.00
19) Phenanthrene-d10	17.116	188	117263	0.400	ng	#-0.01
29) Chrysene-d12	21.314	240	92732	0.400	ng	# 0.00
35) Perylene-d12	23.620	264	64762	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.403	112	63	0.001	ng	0.04
5) Phenol-d6	6.952	99	251	0.005	ng	0.00
8) Nitrobenzene-d5	8.913	82	13559	0.356	ng	0.00
11) 2-Methylnaphthalene-d10	12.138	152	42721	0.273	ng	0.00
14) 2,4,6-Tribromophenol	15.869	330	53	0.126	ng	-0.01
15) 2-Fluorobiphenyl	13.014	172	55241	0.249	ng	0.00
27) Fluoranthene-d10	19.154	212	107325	0.293	ng	0.00
31) Terphenyl-d14	19.755	244	73538	0.362	ng	0.00
Target Compounds						Qvalue
9) Naphthalene	10.598	128	73312	0.324	ng	100
12) 2-Methylnaphthalene	12.213	142	35430	0.237	ng	98
16) Acenaphthylene	14.105	152	142789	0.641	ng	98
17) Acenaphthene	14.448	154	91319	0.593	ng	99
18) Fluorene	15.425	166	160826m	0.860	ng	
24) Pentachlorophenol	16.788	266	453m	0.203	ng	
25) Phenanthrene	17.165	178	2882696	8.963	ng	99
26) Anthracene	17.250	178	591455	2.140	ng	# 94
28) Fluoranthene	19.189	202	4709550	12.940	ng	97
30) Pyrene	19.551	202	4071799	13.602	ng	# 94
32) Benzo(a)anthracene	21.293	228	2127574	7.554	ng	97
33) Chrysene	21.346	228	2019919	6.592	ng	96
34) Bis(2-ethylhexyl)phtha...	21.240	149	1243045	18.606	ng	97
36) Indeno(1,2,3-cd)pyrene	25.988	276	450028	2.140	ng	# 92
37) Benzo(b)fluoranthene	22.925	252	1961583	9.134	ng	# 94
38) Benzo(k)fluoranthene	22.966	252	688874m	3.129	ng	
39) Benzo(a)pyrene	23.517	252	1116465	5.807	ng	# 95
40) Dibenzo(a,h)anthracene	25.995	278	117388	0.683	ng	# 76
41) Benzo(g,h,i)perylene	26.710	276	378934	2.143	ng	95

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

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