

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN113021\
 Data File : BN017642.D
 Acq On : 01 Dec 2021 09:46
 Operator : CG/JU
 Sample : M4862-03
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CS-2B

Manual Integrations
 APPROVED

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

Quant Time: Dec 01 10:31:27 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	25628	0.400	ng	0.00
7) Naphthalene-d8	10.547	136	108835	0.400	ng	# 0.00
13) Acenaphthene-d10	14.384	164	67950	0.400	ng	0.00
19) Phenanthrene-d10	17.116	188	135541	0.400	ng	#-0.01
29) Chrysene-d12	21.303	240	117016	0.400	ng	#-0.01
35) Perylene-d12	23.616	264	110407	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.394	112	6269	0.108	ng	0.03
5) Phenol-d6	6.952	99	13430	0.218	ng	0.00
8) Nitrobenzene-d5	8.913	82	17605	0.381	ng	0.00
11) 2-Methylnaphthalene-d10	12.138	152	55689	0.293	ng	0.00
14) 2,4,6-Tribromophenol	15.869	330	876	0.146	ng	-0.01
15) 2-Fluorobiphenyl	13.014	172	70994	0.268	ng	0.00
27) Fluoranthene-d10	19.154	212	123541	0.292	ng	0.00
31) Terphenyl-d14	19.754	244	89776	0.350	ng	0.00
Target Compounds						Qvalue
9) Naphthalene	10.598	128	297989	1.086	ng	100
12) 2-Methylnaphthalene	12.209	142	85679	0.473	ng	98
16) Acenaphthylene	14.105	152	251917	0.951	ng	99
17) Acenaphthene	14.448	154	79047	0.431	ng	98
18) Fluorene	15.425	166	238293	1.070	ng	95
24) Pentachlorophenol	16.776	266	212	0.197	ng	# 83
25) Phenanthrene	17.165	178	4023776	10.824	ng	99
26) Anthracene	17.250	178	582362	1.823	ng	# 89
28) Fluoranthene	19.188	202	5075406	12.064	ng	97
30) Pyrene	19.546	202	4062587	10.755	ng	97
32) Benzo(a)anthracene	21.293	228	2126502	5.983	ng	97
33) Chrysene	21.346	228	2116558	5.474	ng	97
34) Bis(2-ethylhexyl)phtha...	21.239	149	144844	1.718	ng	95
36) Indeno(1,2,3-cd)pyrene	25.985	276	818835	2.284	ng	# 92
37) Benzo(b)fluoranthene	22.921	252	2602090	7.107	ng	98
38) Benzo(k)fluoranthene	22.962	252	922426m	2.457	ng	
39) Benzo(a)pyrene	23.517	252	1449670	4.423	ng	98
40) Dibenzo(a,h)anthracene	25.995	278	213107	0.727	ng	# 89
41) Benzo(g,h,i)perylene	26.714	276	699184	2.320	ng	95

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

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