Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
Data File : BP009407.D
Acq On : 15 Mar 2022 20:12
Operator : CG/JU
Sample : N1918-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDSH-TB-STOCKPILE-A

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 03/16/2022
Supervised By :Jagrut Upadhyay 03/16/2022

Quant Time: Mar 16 01:58:12 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 05 01:55:59 2022
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.811	152	314293	20.000	ng		-0.01
21) Naphthalene-d8	10.605	136	1273158	20.000	ng		-0.01
39) Acenaphthene-d10	14.451	164	817462	20.000	ng		-0.01
64) Phenanthrene-d10	17.216	188	1674460	20.000	ng		0.00
76) Chrysene-d12	21.333	240	1596872	20.000	ng		0.00
86) Perylene-d12	23.751	264	1875968	20.000	ng		0.02
System Monitoring Compounds							
5) 2-Fluorophenol	5.422	112	831740	38.629	ng		0.01
7) Phenol-d6	7.005	99	1419448	51.876	ng		0.02
23) Nitrobenzene-d5	8.963	82	984503	42.591	ng		-0.01
42) 2,4,6-Tribromophenol	15.951	330	366180	37.097	ng		0.00
45) 2-Fluorobiphenyl	13.075	172	2331601	39.224	ng		-0.01
79) Terphenyl-d14	19.792	244	3376674	39.904	ng		0.00
Target Compounds							Qvalue
10) Phenol	7.034	94	61583	2.173	ng		90
31) Naphthalene	10.658	128	1698580	24.091	ng		99
37) 2-Methylnaphthalene	12.269	142	553552	11.050	ng		98
38) 1-Methylnaphthalene	12.493	142	480922	10.424	ng		99
46) 1,1'-Biphenyl	13.287	154	166819	2.499	ng		98
49) Acenaphthylene	14.175	152	277003	3.546	ng	#	94
52) Acenaphthene	14.516	154	1091907	21.812	ng		99
55) Dibenzofuran	14.851	168	1203133	15.866	ng		97
58) Fluorene	15.504	166	1651792	27.815	ng		97
71) Phenanthrene	17.263	178	16350217m	169.904	ng		
72) Anthracene	17.351	178	4010669	41.871	ng		99
73) Carbazole	17.628	167	1750231	20.053	ng		99
75) Fluoranthene	19.239	202	16021113	144.238	ng		93
78) Pyrene	19.598	202	15903836	147.508	ng	#	86
81) Benzo(a)anthracene	21.316	228	9769335	90.260	ng		97
83) Chrysene	21.369	228	8954369	85.210	ng		97
87) Indeno(1,2,3-cd)pyrene	26.274	276	5759328	39.514	ng	#	91
88) Benzo(b)fluoranthene	23.022	252	11389468m	89.920	ng		
89) Benzo(k)fluoranthene	23.063	252	3471384m	28.688	ng		
90) Benzo(a)pyrene	23.651	252	8486318m	70.641	ng		
91) Dibenzo(a,h)anthracene	26.274	278	1590734	12.871	ng		96
92) Benzo(g,h,i)perylene	27.045	276	5487862	44.947	ng	#	94

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

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