

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
 Data File : BP009422.D
 Acq On : 16 Mar 2022 14:57
 Operator : CG/JU
 Sample : N1918-10DL 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

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

Quant Time: Mar 16 15:58:50 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

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 03/17/2022
 Supervised By : Jagrut Upadhyay 03/17/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.811	152	368173	20.000	ng	-0.01
21) Naphthalene-d8	10.604	136	1318983	20.000	ng	-0.01
39) Acenaphthene-d10	14.451	164	781288	20.000	ng	-0.01
64) Phenanthrene-d10	17.216	188	1597198	20.000	ng	0.00
76) Chrysene-d12	21.327	240	1821998	20.000	ng	0.00
86) Perylene-d12	23.739	264	2092094	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.417	112	904843	35.874	ng	0.00
7) Phenol-d6	7.005	99	1060982	33.101	ng	0.02
23) Nitrobenzene-d5	8.963	82	686321	28.660	ng	-0.01
42) 2,4,6-Tribromophenol	15.951	330	479133	50.787	ng	0.00
45) 2-Fluorobiphenyl	13.075	172	1520219	26.759	ng	-0.01
79) Terphenyl-d14	19.792	244	2349475	24.335	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.652	128	149676	2.049	ng	99
52) Acenaphthene	14.516	154	223336	4.668	ng	100
55) Dibenzofuran	14.851	168	255838	3.530	ng	96
58) Fluorene	15.504	166	334153	5.887	ng	98
71) Phenanthrene	17.257	178	4583419	49.933	ng	99
72) Anthracene	17.345	178	1033415	11.311	ng	99
73) Carbazole	17.622	167	557487	6.696	ng	99
75) Fluoranthene	19.239	202	6213515	58.646	ng	98
78) Pyrene	19.592	202	5669007	46.083	ng	99
81) Benzo(a)anthracene	21.310	228	3485781	28.226	ng	99
83) Chrysene	21.363	228	3121666	26.035	ng	98
87) Indeno(1,2,3-cd)pyrene	26.245	276	2120004	13.043	ng	# 92
88) Benzo(b)fluoranthene	23.010	252	4209097m	29.798	ng	
89) Benzo(k)fluoranthene	23.045	252	1350715m	10.009	ng	
90) Benzo(a)pyrene	23.633	252	3083910	23.019	ng	96
91) Dibenzo(a,h)anthracene	26.262	278	599086	4.347	ng	94
92) Benzo(g,h,i)perylene	27.027	276	2059843	15.128	ng	# 95

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

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