

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021122\
 Data File : BP009129.D
 Acq On : 12 Feb 2022 00:05
 Operator : CG/JU
 Sample : N1505-03 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SU-03-021022

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 02/14/2022
 Supervised By :mohammad ahmed 02/15/2022

Quant Time: Feb 12 00:41:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 07 15:13:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	288892	20.000	ng	-0.01
21) Naphthalene-d8	10.593	136	1191585	20.000	ng	-0.01
39) Acenaphthene-d10	14.439	164	838463	20.000	ng	0.00
64) Phenanthrene-d10	17.198	188	1770227	20.000	ng	0.00
76) Chrysene-d12	21.298	240	1508672	20.000	ng	0.00
86) Perylene-d12	23.698	264	1505667	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	625031	38.593	ng	0.00
7) Phenol-d6	6.993	99	868026	40.676	ng	0.00
23) Nitrobenzene-d5	8.957	82	518106	24.520	ng	-0.01
42) 2,4,6-Tribromophenol	15.939	330	505793	45.180	ng	0.00
45) 2-Fluorobiphenyl	13.057	172	1592735	26.592	ng	-0.01
79) Terphenyl-d14	19.763	244	2411015	28.739	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.640	128	297011	4.887	ng	99
37) 2-Methylnaphthalene	12.257	142	350426	8.148	ng	97
38) 1-Methylnaphthalene	12.475	142	427261	10.167	ng	98
49) Acenaphthylene	14.157	152	836151	11.330	ng	97
52) Acenaphthene	14.504	154	385580	8.587	ng	98
55) Dibenzofuran	14.839	168	359692	4.755	ng	# 87
58) Fluorene	15.492	166	1043040	17.946	ng	99
71) Phenanthrene	17.245	178	5660881	59.058	ng	100
72) Anthracene	17.333	178	1862103	19.856	ng	99
73) Carbazole	17.610	167	213218	2.372	ng	95
75) Fluoranthene	19.221	202	5184476	48.259	ng	97
78) Pyrene	19.574	202	5944357	55.388	ng	99
81) Benzo(a)anthracene	21.280	228	2655520	27.314	ng	98
83) Chrysene	21.333	228	2423724	25.860	ng	98
87) Indeno(1,2,3-cd)pyrene	26.203	276	1185370	10.596	ng	# 93
88) Benzo(b)fluoranthene	22.968	252	2584436m	27.882	ng	
89) Benzo(k)fluoranthene	23.010	252	658723m	7.123	ng	
90) Benzo(a)pyrene	23.598	252	2153333	27.864	ng	98
91) Dibenzo(a,h)anthracene	26.203	278	360929	3.926	ng	# 87
92) Benzo(g,h,i)perylene	26.980	276	1194265	13.047	ng	96

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

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