

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091120\
 Data File : BP003267.D
 Acq On : 11 Sep 2020 18:59
 Operator : CG/JU
 Sample : L3975-04
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BSY-SB-03-0-17

Manual Integrations
 APPROVED

mohammad
 9/14/2020 12:13:29 PM

Quant Time: Sep 11 19:39:07 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 01 14:37:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	226264	20.00	ng	0.00
21) Naphthalene-d8	10.43	136	896897	20.00	ng	0.00
39) Acenaphthene-d10	14.29	164	523766	20.00	ng	0.00
64) Phenanthrene-d10	17.05	188	1055897	20.00	ng	0.00
76) Chrysene-d12	21.14	240	802707	20.00	ng	0.00
86) Perylene-d12	23.37	264	822179	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	976052	76.27	ng	0.00
7) Phenol-d6	6.84	99	1185843	73.92	ng	0.00
23) Nitrobenzene-d5	8.79	82	745900	59.95	ng	0.00
42) 2,4,6-Tribromophenol	15.79	330	576893	81.44	ng	0.00
45) 2-Fluorobiphenyl	12.91	172	1802453	50.40	ng	0.00
79) Terphenyl-d14	19.62	244	1847264	50.64	ng	0.00
Target Compounds						
49) Acenaphthylene	14.01	152	237296	4.880	ng	100
50) Dimethylphthalate	13.75	163	400192	10.019	ng	100
71) Phenanthrene	17.09	178	148671	2.598	ng	99
72) Anthracene	17.18	178	342899	6.262	ng	98
75) Fluoranthene	19.07	202	1669073	25.240	ng	99
78) Pyrene	19.42	202	2024070	38.851	ng	99
81) Benzo(a)anthracene	21.13	228	1098415	21.846	ng	96
83) Chrysene	21.17	228	1214635	25.230	ng	98
87) Indeno(1,2,3-cd)pyrene	25.64	276	817353	13.331	ng	95
88) Benzo(b)fluoranthene	22.71	252	1941007m	37.984	ng	
89) Benzo(k)fluoranthene	22.74	252	674435m	13.814	ng	
90) Benzo(a)pyrene	23.28	252	950941	20.059	ng	98
91) Dibenzo(a,h)anthracene	25.65	278	205766m	4.085	ng	
92) Benzo(g,h,i)perylene	26.34	276	788183	15.593	ng	98

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

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