

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051519\
 Data File : BF114301.D
 Acq On : 15 May 2019 22:45
 Operator : JU/SJ
 Sample : K2766-05DL 4X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P026-SS007-0002-01DL

Manual Integrations
 APPROVED

mohammad
 5/16/2019 9:51:22 AM

Quant Time: May 16 02:04:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 10 15:56:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	243127	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	957624	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	454585	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	846359	20.00	ng	0.00
76) Chrysene-d12	14.02	240	508304	20.00	ng	0.00
87) Perylene-d12	15.50	264	578724	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	571604	37.83	ng	0.00
7) Phenol-d6	6.49	99	786990	44.04	ng	0.00
23) Nitrobenzene-d5	7.41	82	462803	27.12	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	164866	35.08	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	744894	24.79	ng	0.00
79) Terphenyl-d14	12.95	244	557996	22.63	ng	0.00
Target Compounds						
31) Naphthalene	8.15	128	324335	7.251	ng	96
37) 2-Methylnaphthalene	8.85	142	137967	4.780	ng	97
38) 1-Methylnaphthalene	8.95	142	87967	3.237	ng	100
49) Acenaphthylene	9.75	152	493337	10.975	ng	97
50) Dimethylphthalate	9.60	163	123581	3.703	ng	98
58) Fluorene	10.43	166	74512	2.385	ng	99
71) Phenanthrene	11.40	178	1885729	45.796	ng	98
72) Anthracene	11.45	178	319198	7.718	ng	98
75) Fluoranthene	12.60	202	2144654	48.366	ng	100
78) Pyrene	12.83	202	3503368	85.677	ng	99
81) Benzo(a)anthracene	14.01	228	1339603m	40.746	ng	
83) Chrysene	14.05	228	1452476	45.068	ng	98
86) Indeno(1,2,3-cd)pyrene	16.98	276	581738	20.424	ng	97
88) Benzo(b)fluoranthene	15.07	252	1522569m	41.066	ng	
89) Benzo(k)fluoranthene	15.09	252	472615m	13.877	ng	
90) Benzo(a)pyrene	15.44	252	1043280	31.973	ng	100
91) Dibenzo(a,h)anthracene	16.98	278	162153m	5.980	ng	
92) Benzo(g,h,i)perylene	17.43	276	638275	23.490	ng	99

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

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