

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051319\
 Data File : BF114248.D
 Acq On : 14 May 2019 00:33
 Operator : JU/SJ
 Sample : K2766-05DL 2X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 5/14/2019 8:03:35 PM

Quant Time: May 14 13:28:54 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	243623	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	980074	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	456178	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	847256	20.00	ng	0.01
76) Chrysene-d12	14.03	240	543308	20.00	ng	0.00
87) Perylene-d12	15.52	264	578028	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	702489	46.40	ng	0.00
7) Phenol-d6	6.50	99	938728	52.43	ng	0.00
23) Nitrobenzene-d5	7.42	82	549963	31.49	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	189947	40.28	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	850882	28.21	ng	0.00
79) Terphenyl-d14	12.96	244	626147	23.76	ng	0.00
Target Compounds						
31) Naphthalene	8.16	128	382857	8.363	ng	97
37) 2-Methylnaphthalene	8.85	142	164933	5.584	ng	98
38) 1-Methylnaphthalene	8.95	142	108351	3.895	ng	99
49) Acenaphthylene	9.75	152	603012	13.368	ng	98
50) Dimethylphthalate	9.60	163	142722	4.262	ng	98
58) Fluorene	10.44	166	95259	3.038	ng	93
71) Phenanthrene	11.40	178	2102866	51.016	ng	99
72) Anthracene	11.45	178	382147	9.230	ng	98
75) Fluoranthene	12.60	202	2438475	54.935	ng	97
78) Pyrene	12.84	202	3911495	89.495	ng	97
81) Benzo(a)anthracene	14.02	228	1531736m	43.588	ng	
83) Chrysene	14.06	228	1758800	51.057	ng	99
86) Indeno(1,2,3-cd)pyrene	17.00	276	660451	21.694	ng	97
88) Benzo(b)fluoranthene	15.09	252	1719128m	46.423	ng	
89) Benzo(k)fluoranthene	15.10	252	545107m	16.024	ng	
90) Benzo(a)pyrene	15.46	252	1158579	35.549	ng	99
91) Dibenzo(a,h)anthracene	17.00	278	180254m	6.656	ng	
92) Benzo(g,h,i)perylene	17.46	276	747415	27.539	ng	99

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

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