

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051219\
 Data File : BF114214.D
 Acq On : 13 May 2019 1:40
 Operator : JU/SJ
 Sample : K2764-16DL 2X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P026-SS002-0002-01DL

Manual Integrations
 APPROVED

Sohil
 5/13/2019 8:37:48 PM

Quant Time: May 13 09:28:13 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	272943	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	1041925	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	474322	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	837676	20.00	ng	0.00
76) Chrysene-d12	14.02	240	593197	20.00	ng	0.00
87) Perylene-d12	15.49	264	622731	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	790646	46.62	ng	0.00
7) Phenol-d6	6.49	99	1052791	52.48	ng	0.00
23) Nitrobenzene-d5	7.41	82	639106	34.42	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	211102	43.05	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	947782	30.23	ng	0.00
79) Terphenyl-d14	12.95	244	715263	24.86	ng	0.00
Target Compounds						
31) Naphthalene	8.15	128	287959	5.917	ng	97
37) 2-Methylnaphthalene	8.85	142	111330	3.545	ng	96
38) 1-Methylnaphthalene	8.95	142	71021	2.402	ng	99
49) Acenaphthylene	9.75	152	536109	11.430	ng	97
50) Dimethylphthalate	9.60	163	105965	3.043	ng	99
58) Fluorene	10.43	166	67980	2.085	ng	98
71) Phenanthrene	11.40	178	2021968	49.614	ng	99
72) Anthracene	11.45	178	335131	8.187	ng	99
75) Fluoranthene	12.59	202	2593331	59.091	ng	96
78) Pyrene	12.83	202	4237601	88.802	ng	96
81) Benzo(a)anthracene	14.01	228	1680904m	43.810	ng	
83) Chrysene	14.04	228	1979448	52.629	ng	99
86) Indeno(1,2,3-cd)pyrene	16.96	276	753777	22.677	ng	98
88) Benzo(b)fluoranthene	15.07	252	1829153m	45.848	ng	
89) Benzo(k)fluoranthene	15.08	252	644143m	17.576	ng	
90) Benzo(a)pyrene	15.43	252	1185459	33.763	ng	100
91) Dibenzo(a,h)anthracene	16.97	278	210159	7.203	ng	97
92) Benzo(g,h,i)perylene	17.42	276	867547	29.671	ng	98

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

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