

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060519\
 Data File : BF114846.D
 Acq On : 6 Jun 2019 3:56
 Operator : HP/JU
 Sample : K3151-01 2X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-060419-A

Manual Integrations
 APPROVED

mohammad
 6/6/2019 4:50:56 PM

Quant Time: Jun 06 08:09:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF060419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 04 16:57:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	381144	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	1448774	20.00	ng	0.00
39) Acenaphthene-d10	9.70	164	681054	20.00	ng	0.00
64) Phenanthrene-d10	11.17	188	1207857	20.00	ng	0.00
76) Chrysene-d12	13.80	240	793190	20.00	ng	0.00
87) Perylene-d12	15.19	264	619240	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	1118361	51.27	ng	0.00
7) Phenol-d6	6.30	99	1382866	52.10	ng	0.00
23) Nitrobenzene-d5	7.23	82	804203	34.17	ng	0.00
42) 2,4,6-Tribromophenol	10.48	330	341451	46.52	ng	0.00
45) 2-Fluorobiphenyl	9.03	172	1556268	42.81	ng	0.00
79) Terphenyl-d14	12.76	244	1458576	35.76	ng	0.01
Target Compounds						
50) Dimethylphthalate	9.42	163	274621	5.521	ng	98
71) Phenanthrene	11.20	178	668207	11.588	ng	96
72) Anthracene	11.25	178	190175	3.134	ng	94
75) Fluoranthene	12.38	202	913574	15.222	ng	90
78) Pyrene	12.61	202	1118870	16.260	ng	96
81) Benzo(a)anthracene	13.79	228	545074	8.918	ng	98
83) Chrysene	13.83	228	549382	9.239	ng	96
86) Indeno(1,2,3-cd)pyrene	16.50	276	252609	3.723	ng	94
88) Benzo(b)fluoranthene	14.80	252	542731m	13.767	ng	
89) Benzo(k)fluoranthene	14.82	252	203599m	6.082	ng	
90) Benzo(a)pyrene	15.13	252	390409	11.442	ng	97
91) Dibenzo(a,h)anthracene	16.50	278	67307	2.083	ng	# 73
92) Benzo(a,h,i)perylene	16.89	276	233939	7.199	ng	97

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

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