

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
 Data File : BF114349.D
 Acq On : 17 May 2019 15:01
 Operator : JU/SJ
 Sample : K2644-03 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-051519-B

Manual Integrations
 APPROVED

mohammad
 5/21/2019 2:15:48 PM

Quant Time: May 18 05:17:08 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	358934	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	1396910	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	663593	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	1309789	20.00	ng	0.00
76) Chrysene-d12	14.02	240	896726	20.00	ng	0.00
87) Perylene-d12	15.51	264	1038153	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1116581	50.06	ng	0.00
7) Phenol-d6	6.49	99	1458685	55.30	ng	0.00
23) Nitrobenzene-d5	7.40	82	883335	35.49	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	355564	51.83	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1462757	33.34	ng	-0.01
79) Terphenyl-d14	12.95	244	1401653	32.22	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	261712	5.372	ng	98
71) Phenanthrene	11.39	178	618287	9.703	ng	98
72) Anthracene	11.44	178	180758	2.824	ng	98
73) Carbazole	11.60	167	137416	2.251	ng	96
75) Fluoranthene	12.58	202	781902	11.394	ng	99
78) Pyrene	12.81	202	608746	8.439	ng	98
81) Benzo(a)anthracene	14.00	228	295070	5.087	ng	99
83) Chrysene	14.04	228	263938	4.642	ng	98
86) Indeno(1,2,3-cd)pyrene	16.99	276	148986	2.965	ng	99
88) Benzo(b)fluoranthene	15.07	252	381536m	5.737	ng	
90) Benzo(a)pyrene	15.44	252	261414	4.466	ng	99
92) Benzo(g,h,i)perylene	17.43	276	147994	3.036	ng	96

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

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