

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
 Data File : BF114373.D
 Acq On : 18 May 2019 2:29
 Operator : JU/SJ
 Sample : K2894-05
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-01-0-1

Manual Integrations
 APPROVED

mohammad
 5/21/2019 2:16:40 PM

Quant Time: May 18 06:37:26 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	252747	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	943638	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	407885	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	672415	20.00	ng	0.00
76) Chrysene-d12	14.02	240	418819	20.00	ng	0.00
87) Perylene-d12	15.49	264	316015	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	1287089	81.95	ng	0.00
7) Phenol-d6	6.50	99	1674061	90.12	ng	0.00
23) Nitrobenzene-d5	7.41	82	597570	35.54	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	188281	44.65	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1628763	60.40	ng	0.00
79) Terphenyl-d14	12.96	244	1081184	53.22	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.60	163	256216	8.557	ng	99
71) Phenanthrene	11.40	178	179780	5.496	ng	98
72) Anthracene	11.45	178	67943	2.068	ng	98
75) Fluoranthene	12.59	202	370859	10.527	ng	97
78) Pyrene	12.82	202	348566	10.346	ng	98
81) Benzo(a)anthracene	14.00	228	170915	6.309	ng	98
83) Chrysene	14.04	228	161656	6.088	ng	99
86) Indeno(1,2,3-cd)pyrene	16.98	276	63350	2.699	ng	98
88) Benzo(b)fluoranthene	15.06	252	175733m	8.680	ng	
89) Benzo(k)fluoranthene	15.09	252	58917m	3.168	ng	
90) Benzo(a)pyrene	15.43	252	136252	7.647	ng	# 78
92) Benzo(g,h,i)perylene	17.43	276	54183	3.652	ng	# 79

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

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