

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051719\
 Data File : BF114374.D
 Acq On : 18 May 2019 2:55
 Operator : JU/SJ
 Sample : K2894-06
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-02-0-1

Manual Integrations
 APPROVED

mohammad
 5/21/2019 2:16:43 PM

Quant Time: May 18 06:39:33 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	215491	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	841986	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	382954	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	702244	20.00	ng	0.00
76) Chrysene-d12	14.02	240	483317	20.00	ng	0.00
87) Perylene-d12	15.49	264	391779	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	805232	60.13	ng	0.00
7) Phenol-d6	6.49	99	1053469	66.52	ng	0.00
23) Nitrobenzene-d5	7.40	82	329310	21.95	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	125072	31.59	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1217032	48.07	ng	0.00
79) Terphenyl-d14	12.96	244	1107586	47.24	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	148077	5.267	ng	99
71) Phenanthrene	11.40	178	304321	8.907	ng	97
72) Anthracene	11.45	178	109235	3.183	ng	96
75) Fluoranthene	12.59	202	683153	18.568	ng	98
78) Pyrene	12.82	202	601790	15.478	ng	97
81) Benzo(a)anthracene	14.00	228	346505	11.084	ng	99
83) Chrysene	14.05	228	297093	9.695	ng	98
86) Indeno(1,2,3-cd)pyrene	16.97	276	106417	3.929	ng	98
88) Benzo(b)fluoranthene	15.06	252	324970m	12.947	ng	
89) Benzo(k)fluoranthene	15.09	252	100163m	4.344	ng	
90) Benzo(a)pyrene	15.43	252	225471	10.207	ng	# 93
92) Benzo(g,h,i)perylene	17.43	276	93667	5.092	ng	# 94

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

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