

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052219\
 Data File : BF114506.D
 Acq On : 23 May 2019 1:39
 Operator : JU/SJ
 Sample : K2970-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-11-S

Manual Integrations
 APPROVED

mohammad
 5/24/2019 8:49:28 AM

Quant Time: May 23 04:39:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 20 04:47:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	184867	20.00	ng	-0.01
21) Naphthalene-d8	8.11	136	720672	20.00	ng	-0.01
39) Acenaphthene-d10	9.86	164	335095	20.00	ng	-0.01
64) Phenanthrene-d10	11.35	188	617337	20.00	ng	0.00
76) Chrysene-d12	14.00	240	408672	20.00	ng	0.00
87) Perylene-d12	15.46	264	423596	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	914996	79.65	ng	0.00
7) Phenol-d6	6.47	99	1190720	87.64	ng	-0.01
23) Nitrobenzene-d5	7.39	82	687338	53.52	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	252369	72.85	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	978489	44.17	ng	0.00
79) Terphenyl-d14	12.93	244	737631	37.21	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.58	163	173042	7.034	ng	99
71) Phenanthrene	11.37	178	112329	3.740	ng	97
75) Fluoranthene	12.57	202	250816	7.755	ng	98
78) Pyrene	12.80	202	254856	7.752	ng	98
81) Benzo(a)anthracene	13.99	228	125582	4.751	ng	98
83) Chrysene	14.02	228	108676	4.194	ng	98
88) Benzo(b)fluoranthene	15.04	252	143513m	5.288	ng	
90) Benzo(a)pyrene	15.40	252	103634	4.339	ng	97
92) Benzo(g,h,i)perylene	17.40	276	47408	2.384	ng	# 92

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

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