

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021219\
 Data File : BF112516.D
 Acq On : 12 Feb 2019 19:30
 Operator : JU/SJ
 Sample : K1343-07 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-021119-D

Manual Integrations
 APPROVED

Sohil
 2/13/2019 11:12:17 AM

Quant Time: Feb 13 04:43:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	120153	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	461987	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	204004	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	330409	20.00	ng	0.00
76) Chrysene-d12	14.01	240	278214	20.00	ng	0.00
87) Perylene-d12	15.49	264	232570	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	385007	68.06	ng	-0.01
7) Phenol-d6	6.48	99	479828	61.04	ng	-0.01
23) Nitrobenzene-d5	7.40	82	298336	37.21	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	113759	47.91	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	563058	39.04	ng	-0.01
79) Terphenyl-d14	12.94	244	463895	31.40	ng	0.00
Target Compounds						
10) Phenol	6.50	94	24938	2.892	ng	94
50) Dimethylphthalate	9.59	163	64940	4.097	ng	97
71) Phenanthrene	11.39	178	191891	11.171	ng	97
72) Anthracene	11.44	178	60905	3.559	ng	97
75) Fluoranthene	12.58	202	261133	14.174	ng	97
78) Pyrene	12.81	202	219334	11.387	ng	98
81) Benzo(a)anthracene	14.00	228	120149	7.346	ng	99
83) Chrysene	14.03	228	108776	6.968	ng	97
86) Indeno(1,2,3-cd)pyrene	16.99	276	29792	2.408	ng	98
88) Benzo(b)fluoranthene	15.06	252	101393m	7.290	ng	
89) Benzo(k)fluoranthene	15.08	252	36447m	2.736	ng	
90) Benzo(a)pyrene	15.43	252	70086	5.600	ng	# 91
92) Benzo(a,h,i)perylene	17.45	276	29875	3.139	ng	# 87

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

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