

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032819\
 Data File : BF113396.D
 Acq On : 29 Mar 2019 3:14
 Operator : JU/SJ
 Sample : K2086-07DL 2X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-18DL

Manual Integrations
 APPROVED

Sohil
 3/29/2019 4:47:27 PM

Quant Time: Mar 29 05:33:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	175037	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	682155	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	326541	20.00	ng	0.00
64) Phenanthrene-d10	11.21	188	616895	20.00	ng	0.00
76) Chrysene-d12	13.84	240	431883	20.00	ng	0.00
87) Perylene-d12	15.24	264	464908	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	566989	52.97	ng	0.00
7) Phenol-d6	6.35	99	740331	54.67	ng	-0.01
23) Nitrobenzene-d5	7.26	82	430275	37.44	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	215604	53.45	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	898175	32.14	ng	0.00
79) Terphenyl-d14	12.79	244	823585	42.03	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.45	163	56765	2.346	ng	99
71) Phenanthrene	11.23	178	607822	19.841	ng	99
72) Anthracene	11.29	178	89416	2.874	ng	97
73) Carbazole	11.45	167	125529	4.228	ng	98
75) Fluoranthene	12.42	202	1733947	50.067	ng	98
78) Pyrene	12.64	202	1374618	45.931	ng	99
81) Benzo(a)anthracene	13.83	228	262671	10.625	ng	95
83) Chrysene	13.87	228	927537	36.938	ng	99
86) Indeno(1,2,3-cd)pyrene	16.59	276	217834	10.108	ng	98
88) Benzo(b)fluoranthene	14.85	252	1110595m	38.456	ng	
89) Benzo(k)fluoranthene	14.87	252	254056m	10.017	ng	
90) Benzo(a)pyrene	15.19	252	231588	9.045	ng	97
91) Dibenzo(a,h)anthracene	16.60	278	48528	2.192	ng	# 80
92) Benzo(g,h,i)perylene	17.00	276	181028	8.110	ng	98

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

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