

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
 Data File : BF108088.D
 Acq On : 10 Aug 2018 17:10
 Operator : JU/SJ
 Sample : J4272-05 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-03-080918-C

Manual Integrations
 APPROVED

Sohil
 8/13/2018 3:51:59 PM

Quant Time: Aug 10 18:18:04 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 12:24:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.02	152	283465	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	1013867	20.00	ng	0.00
38) Acenaphthene-d10	10.07	164	457508	20.00	ng	0.00
63) Phenanthrene-d10	11.56	188	753532	20.00	ng	0.00
75) Chrysene-d12	14.22	240	671586	20.00	ng	0.00
86) Perylene-d12	15.79	264	586316	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	730816	46.68	ng	0.00
7) Phenol-d6	6.62	99	968961	46.57	ng	0.00
23) Nitrobenzene-d5	7.57	82	612978	33.74	ng	0.00
41) 2,4,6-Tribromophenol	10.85	330	191789	42.03	ng	0.00
44) 2-Fluorobiphenyl	9.37	172	947846	33.26	ng	0.00
78) Terphenyl-d14	13.15	244	833265	31.55	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.77	163	95459	2.995	ng	Qvalue # 99
51) Acenaphthene	10.10	154	57677	2.234	ng	99
57) Fluorene	10.61	166	112818	3.810	ng	100
70) Phenanthrene	11.58	178	898310	25.275	ng	99
71) Anthracene	11.64	178	270831	7.435	ng	99
74) Fluoranthene	12.78	202	1207246	31.123	ng	96
77) Pyrene	13.02	202	1131699	26.617	ng	99
80) Benzo(a)anthracene	14.21	228	602296	15.627	ng	97
82) Chrysene	14.24	228	500066	14.152	ng	97
85) Indeno(1,2,3-cd)pyrene	17.41	276	166545	5.440	ng	# 89
87) Benzo(b)fluoranthene	15.31	252	608800m	17.377	ng	
88) Benzo(k)fluoranthene	15.34	252	156760m	4.867	ng	
89) Benzo(a)pyrene	15.72	252	389509	12.416	ng	98
91) Benzo(g,h,i)perylene	17.90	276	150132	5.874	ng	# 93

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

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