

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081318\
 Data File : BF108124.D
 Acq On : 13 Aug 2018 18:53
 Operator : JU/SJ
 Sample : J4430-12
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G.L-WC-01

Manual Integrations
 APPROVED

Sohil
 8/14/2018 2:36:49 PM

Quant Time: Aug 14 01:46:30 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	206216	20.00	ng	0.01
21) Naphthalene-d8	8.31	136	733334	20.00	ng	0.00
38) Acenaphthene-d10	10.07	164	341072	20.00	ng	0.00
63) Phenanthrene-d10	11.57	188	602279	20.00	ng	0.01
75) Chrysene-d12	14.24	240	476842	20.00	ng	0.02
86) Perylene-d12	15.81	264	334356	20.00	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.65	112	1279750	112.35	ng	0.02
7) Phenol-d6	6.64	99	1702408	112.47	ng	0.01
23) Nitrobenzene-d5	7.58	82	1073212	81.66	ng	0.00
41) 2,4,6-Tribromophenol	10.87	330	365704	107.49	ng	0.01
44) 2-Fluorobiphenyl	9.38	172	1669848	78.60	ng	0.00
78) Terphenyl-d14	13.17	244	1527183	81.43	ng	0.02
Target Compounds						
49) Dimethylphthalate	9.77	163	233490	9.828	ng	99
70) Phenanthrene	11.60	178	287058	10.105	ng	99
71) Anthracene	11.65	178	69106	2.374	ng	99
74) Fluoranthene	12.80	202	381276	12.298	ng	96
77) Pyrene	13.03	202	384795	12.746	ng	99
80) Benzo(a)anthracene	14.22	228	166876	6.098	ng	98
82) Chrysene	14.26	228	153998	6.138	ng	97
85) Indeno(1,2,3-cd)pyrene	17.45	276	44412	2.043	ng	# 85
87) Benzo(b)fluoranthene	15.34	252	124566m	6.235	ng	
88) Benzo(k)fluoranthene	15.36	252	40148m	2.186	ng	
89) Benzo(a)pyrene	15.74	252	85997	4.807	ng	97
91) Benzo(g,h,i)perylene	17.95	276	43922	3.013	ng	# 90

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

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