

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041919\
 Data File : BF113862.D
 Acq On : 19 Apr 2019 16:55
 Operator : JU/SJ
 Sample : K2202-03 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SA-06-041819-B

Manual Integrations
 APPROVED

Sohil
 4/22/2019 7:26:51 PM

Quant Time: Apr 20 04:40:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 13:55:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	194141	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	671877	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	302102	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	550756	20.00	ng	0.02
76) Chrysene-d12	14.20	240	501355	20.00	ng	0.16
87) Perylene-d12	15.94	264	392808	20.00	ng	0.44
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	586837	50.26	ng	0.01
7) Phenol-d6	6.52	99	749474	51.47	ng	0.00
23) Nitrobenzene-d5	7.46	82	406394	36.83	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	174532	52.29	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	696445	35.16	ng	0.00
79) Terphenyl-d14	13.04	244	729778	30.52	ng	0.04
Target Compounds						
50) Dimethylphthalate	9.65	163	77740	3.677	ng	Qvalue 100
71) Phenanthrene	11.45	178	84928	3.028	ng	98
75) Fluoranthene	12.65	202	160482	5.398	ng	96
78) Pyrene	12.89	202	177239	5.016	ng	97
81) Benzo(a)anthracene	14.19	228	94753	3.233	ng	94
83) Chrysene	14.23	228	83192	2.860	ng	98
88) Benzo(b)fluoranthene	15.47	252	94233m	3.848	ng	
90) Benzo(a)pyrene	15.87	252	70429m	3.253	ng	
92) Benzo(g,h,i)perylene	17.94	276	41659	2.127	ng	# 90

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

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