

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041819\
 Data File : BF113839.D
 Acq On : 19 Apr 2019 1:48
 Operator : JU/SJ
 Sample : K2197-11 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-041619-D

Manual Integrations
 APPROVED

Sohil
 4/23/2019 8:16:54 AM

Quant Time: Apr 19 04:59:38 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	148564	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	501366	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	241062	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	524963	20.00	ng	0.00
76) Chrysene-d12	14.04	240	521099	20.00	ng	0.00
87) Perylene-d12	15.52	264	401512	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	413286	46.26	ng	0.00
7) Phenol-d6	6.51	99	496898	44.60	ng	0.00
23) Nitrobenzene-d5	7.45	82	264232	32.09	ng	-0.01
42) 2,4,6-Tribromophenol	10.71	330	142898	53.65	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	515441	32.61	ng	0.00
79) Terphenyl-d14	12.99	244	773017	31.11	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.64	163	57573	3.413	ng	98
71) Phenanthrene	11.43	178	83571	3.126	ng	98
75) Fluoranthene	12.62	202	178622	6.304	ng	98
78) Pyrene	12.85	202	158090	4.305	ng	97
81) Benzo(a)anthracene	14.03	228	91908	3.017	ng	97
83) Chrysene	14.07	228	80447	2.661	ng	96
88) Benzo(b)fluoranthene	15.09	252	77068m	3.079	ng	
90) Benzo(a)pyrene	15.45	252	48444	2.189	ng	# 93

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

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