

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062719\
 Data File : BF115258.D
 Acq On : 28 Jun 2019 2:14
 Operator : HP/JU
 Sample : K3157-15 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-062619-C

Manual Integrations
 APPROVED

mohammad
 6/28/2019 2:59:01 PM

Quant Time: Jun 28 07:43:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 28 05:26:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	240565	20.00	ng	0.00
21) Naphthalene-d8	7.88	136	937555	20.00	ng	0.00
39) Acenaphthene-d10	9.62	164	472821	20.00	ng	0.00
64) Phenanthrene-d10	11.10	188	942042	20.00	ng	0.00
76) Chrysene-d12	13.72	240	697512	20.00	ng	0.00
87) Perylene-d12	15.08	264	560603	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	668225	42.87	ng	0.00
7) Phenol-d6	6.24	99	862697	45.59	ng	0.00
23) Nitrobenzene-d5	7.15	82	495650	32.52	ng	-0.01
42) 2,4,6-Tribromophenol	10.41	330	259597	52.06	ng	0.00
45) 2-Fluorobiphenyl	8.95	172	994169	35.72	ng	0.00
79) Terphenyl-d14	12.68	244	1145347	34.91	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.35	163	194277	5.585	ng	99
71) Phenanthrene	11.12	178	274040	5.403	ng	96
75) Fluoranthene	12.30	202	375940	7.429	ng	98
78) Pyrene	12.52	202	387527	7.229	ng	98
81) Benzo(a)anthracene	13.71	228	167113	3.339	ng	93
83) Chrysene	13.74	228	170848	3.727	ng	95
88) Benzo(b)fluoranthene	14.71	252	177869m	5.085	ng	
90) Benzo(a)pyrene	15.02	252	120300	3.816	ng	# 89

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

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