

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061319\
 Data File : BF115025.D
 Acq On : 14 Jun 2019 1:50
 Operator : HP/JU
 Sample : K3156-03 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HD-01-061219

Manual Integrations
 APPROVED

mohammad
 6/14/2019 8:47:31 AM

Quant Time: Jun 14 05:12:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 12 15:42:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	139576	20.00	ng	0.00
21) Naphthalene-d8	7.93	136	522205	20.00	ng	0.00
39) Acenaphthene-d10	9.67	164	244227	20.00	ng	0.00
64) Phenanthrene-d10	11.15	188	403272	20.00	ng	-0.01
76) Chrysene-d12	13.77	240	253350	20.00	ng	-0.01
87) Perylene-d12	15.14	264	239421	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	909182	108.77	ng	0.00
7) Phenol-d6	6.29	99	1129159	111.98	ng	-0.01
23) Nitrobenzene-d5	7.20	82	680470	78.43	ng	-0.01
42) 2,4,6-Tribromophenol	10.46	330	240731	92.05	ng	-0.01
45) 2-Fluorobiphenyl	9.00	172	1256300	87.18	ng	0.00
79) Terphenyl-d14	12.73	244	945194	70.17	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.40	163	267808	14.581	ng	97
71) Phenanthrene	11.17	178	53077	2.549	ng	95
75) Fluoranthene	12.35	202	111775	5.250	ng	97
78) Pyrene	12.57	202	95841	4.153	ng	97
81) Benzo(a)anthracene	13.76	228	53451	2.819	ng	94
83) Chrysene	13.79	228	62105	3.294	ng	95
88) Benzo(b)fluoranthene	14.76	252	84581m	5.314	ng	
90) Benzo(a)pyrene	15.09	252	50153	3.652	ng	# 94
92) Benzo(g,h,i)perylene	16.83	276	29101	2.413	ng	97

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

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