

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060719\
 Data File : BF114916.D
 Acq On : 8 Jun 2019 3:11
 Operator : HP/JU
 Sample : K3250-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CHIMNEY-1

Quant Time: Jun 10 08:36:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF060419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 04 16:57:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	306576	20.00	ng	0.00
21) Naphthalene-d8	7.94	136	1133327	20.00	ng	-0.01
39) Acenaphthene-d10	9.69	164	506049	20.00	ng	-0.01
64) Phenanthrene-d10	11.16	188	809721	20.00	ng	0.00
76) Chrysene-d12	13.79	240	588442	20.00	ng	0.00
87) Perylene-d12	15.17	264	438696	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	1000008	57.00	ng	0.01
7) Phenol-d6	6.30	99	1217291	57.02	ng	0.00
23) Nitrobenzene-d5	7.22	82	700096	38.03	ng	-0.01
42) 2,4,6-Tribromophenol	10.47	330	256678	47.07	ng	0.00
45) 2-Fluorobiphenyl	9.02	172	1337926	49.53	ng	-0.01
79) Terphenyl-d14	12.75	244	1082123	35.76	ng	0.00
Target Compounds						
32) Benzoic acid	7.73	122	211646	18.508	ng	98
37) 2-Methylnaphthalene	8.65	142	131943	3.728	ng	98
50) Dimethylphthalate	9.41	163	84976	2.299	ng	98
56) 4-Nitrophenol	9.82	139	98842	13.284	ng	94
71) Phenanthrene	11.19	178	293746	7.599	ng	99
75) Fluoranthene	12.37	202	160829	3.997	ng	95
78) Pyrene	12.60	202	146466	2.869	ng	97
83) Chrysene	13.81	228	95697	2.169	ng	95
88) Benzo(b)fluoranthene	14.79	252	91557m	3.278	ng	
92) Benzo(q,h,i)perylene	16.87	276	46502	2.020	ng	# 90

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

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