

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060619\
 Data File : BF114861.D
 Acq On : 6 Jun 2019 14:22
 Operator : HP/JU
 Sample : K3189-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-7A

Manual Integrations
 APPROVED

mohammad
 6/7/2019 2:11:28 PM

Quant Time: Jun 07 01:58:48 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	6.67	152	469019	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	1807516	20.00	ng	0.00
39) Acenaphthene-d10	9.70	164	867134	20.00	ng	0.00
64) Phenanthrene-d10	11.17	188	1505479	20.00	ng	0.00
76) Chrysene-d12	13.79	240	901033	20.00	ng	0.00
87) Perylene-d12	15.19	264	918401	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	3105472	115.70	ng	0.02
7) Phenol-d6	6.32	99	3846695	117.78	ng	0.01
23) Nitrobenzene-d5	7.23	82	2450000	83.44	ng	0.00
42) 2,4,6-Tribromophenol	10.49	330	928796	99.39	ng	0.00
45) 2-Fluorobiphenyl	9.03	172	3388305	73.20	ng	0.00
79) Terphenyl-d14	12.76	244	2900878	62.61	ng	0.00
Target Compounds						
32) Benzoic acid	7.67	122	74796m	4.101	ng	
50) Dimethylphthalate	9.42	163	900183	14.214	ng	99
75) Fluoranthene	12.37	202	323047	4.319	ng	95
78) Pyrene	12.60	202	274449	3.511	ng	96
81) Benzo(a)anthracene	13.78	228	162822	2.345	ng	97
83) Chrysene	13.82	228	143604	2.126	ng	97
88) Benzo(b)fluoranthene	14.80	252	178122m	3.047	ng	
90) Benzo(a)pyrene	15.12	252	123867	2.448	ng	# 95

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

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