

Data Path : Z:\HPCHEM1\BNA F\DATA\BF053117\
 Data File : BF095590.D
 Acq On : 1 Jun 2017 7:01
 Operator : SJ/MA
 Sample : I3388-06
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-B29-B35-002

Manual Integrations
 APPROVED

mohammad
 6/1/2017 2:20:54 PM

Quant Time: Jun 01 08:18:34 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 01 04:09:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	92444	20.00	ng	0.00
21) Naphthalene-d8	9.58	136	299295	20.00	ng	0.00
38) Acenaphthene-d10	12.41	164	124977	20.00	ng	0.00
63) Phenanthrene-d10	14.81	188	192462	20.00	ng	0.00
75) Chrysene-d12	18.51	240	158142	20.00	ng	0.00
86) Perylene-d12	20.18	264	118342	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	705196	127.18	ng	0.03
7) Phenol-d6	7.11	99	806166	121.17	ng	0.00
23) Nitrobenzene-d5	8.47	82	430453	90.69	ng	0.00
41) 2,4,6-Tribromophenol	13.71	330	136351	133.22	ng	0.00
44) 2-Fluorobiphenyl	11.36	172	854938	98.46	ng	0.00
78) Terphenyl-d14	17.16	244	577749	80.47	ng	0.00
Target Compounds						
17) 2-Methylphenol	8.14	107	15236	2.95	ng	Qvalue 89
20) 3+4-Methylphenols	8.42	107	22029	3.14	ng	# 57
27) 2,4-Dimethylphenol	9.17	122	42166m	9.72	ng	
31) Naphthalene	9.62	128	1213157	80.65	ng	99
37) 2-Methylnaphthalene	10.74	142	115356	11.68	ng	97
51) Acenaphthene	12.46	154	28250	3.55	ng	98
57) Fluorene	13.31	166	19693	2.06	ng	# 92
70) Phenanthrene	14.85	178	26595	2.40	ng	99
72) Carbazole	15.24	167	67244	6.54	ng	# 93

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

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