

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090117\
 Data File : BF098221.D
 Acq On : 1 Sep 2017 12:46
 Operator : SJ/JU
 Sample : I4987-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-06-COMP

Manual Integrations
 APPROVED

mohammad
 9/5/2017 5:21:58 PM

Quant Time: Sep 01 16:26:40 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 01 16:25:25 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	112435	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	443992	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	188908	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	308760	20.00	ng	0.00
75) Chrysene-d12	13.87	240	211187	20.00	ng	0.00
86) Perylene-d12	15.28	264	202934	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	892645	120.76	ng	0.00
7) Phenol-d6	6.37	99	1218470	121.32	ng	0.00
23) Nitrobenzene-d5	7.29	82	746416	91.33	ng	0.00
41) 2,4,6-Tribromophenol	10.54	330	192850	117.66	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	900294	70.02	ng	0.00
78) Terphenyl-d14	12.82	244	550463	54.35	ng	0.00
Target Compounds						
10) Phenol	6.38	94	33325	2.837	ng	89
49) Dimethylphthalate	9.47	163	233027	16.875	ng	98
70) Phenanthrene	11.26	178	80472	4.891	ng	99
74) Fluoranthene	12.44	202	108728	6.331	ng	98
77) Pyrene	12.67	202	102594	5.940	ng	98
80) Benzo(a)anthracene	13.86	228	52113	4.117	ng	100
82) Chrysene	13.89	228	45963m	3.650	ng	
83) Bis(2-ethylhexyl)phthalate	13.86	149	41266	5.623	ng	# 99
87) Benzo(b)fluoranthene	14.88	252	52585m	4.111	ng	
89) Benzo(a)pyrene	15.22	252	38784	3.425	ng	98

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

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