

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060617\
 Data File : BF095742.D
 Acq On : 7 Jun 2017 11:29
 Operator : SJ/MA
 Sample : I3469-01
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-2(0-2)20170603

Manual Integrations
 APPROVED

mohammad
 6/7/2017 2:19:36 PM

Quant Time: Jun 07 13:51:45 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 16:15:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	73462	20.00	ng	-0.01
21) Naphthalene-d8	9.41	136	276545	20.00	ng	-0.02
38) Acenaphthene-d10	12.23	164	106908	20.00	ng	-0.02
63) Phenanthrene-d10	14.63	188	161661	20.00	ng	-0.02
75) Chrysene-d12	18.34	240	143465	20.00	ng	-0.01
86) Perylene-d12	20.01	264	99908	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	695412	150.84	ng	0.00
7) Phenol-d6	6.95	99	814433	147.56	ng	0.00
23) Nitrobenzene-d5	8.30	82	454343	102.03	ng	-0.02
41) 2,4,6-Tribromophenol	13.53	330	120775	127.68	ng	-0.02
44) 2-Fluorobiphenyl	11.20	172	803402	104.58	ng	-0.02
78) Terphenyl-d14	17.00	244	518062	89.62	ng	-0.01
Target Compounds						
10) Phenol	6.96	94	12034	2.102	ng	Qvalue # 74
49) Dimethylphthalate	11.89	163	120050	14.790	ng	99
70) Phenanthrene	14.67	178	25161	2.697	ng	96
74) Fluoranthene	16.44	202	74674	8.088	ng	98
77) Pyrene	16.74	202	67579	6.313	ng	96
80) Benzo(a)anthracene	18.33	228	38578	4.625	ng	94
82) Chrysene	18.38	228	35613	4.272	ng	97
85) Indeno(1,2,3-cd)pyrene	21.19	276	21679	3.040	ng	97
87) Benzo(b)fluoranthene	19.61	252	41538m	6.640	ng	
88) Benzo(k)fluoranthene	19.64	252	16754m	2.802	ng	
89) Benzo(a)pyrene	19.96	252	32491	5.651	ng	97
91) Benzo(g,h,i)perylene	21.53	276	21880	4.400	ng	98

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

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