

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092517\
 Data File : BF098805.D
 Acq On : 26 Sep 2017 1:49
 Operator : SJ/JU
 Sample : I5362-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-2-A-COMP

Manual Integrations
 APPROVED

Sohil
 9/26/2017 4:58:50 PM

Quant Time: Sep 26 07:54:35 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 23 13:50:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	148106	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	623705	20.00	ng	-0.01
38) Acenaphthene-d10	9.95	164	279993	20.00	ng	-0.01
63) Phenanthrene-d10	11.44	188	458462	20.00	ng	0.00
75) Chrysene-d12	14.07	240	264839	20.00	ng	-0.01
86) Perylene-d12	15.57	264	260577	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	904938	97.27	ng	0.00
7) Phenol-d6	6.54	99	1107894	96.53	ng	0.00
23) Nitrobenzene-d5	7.47	82	680298	69.95	ng	-0.01
41) 2,4,6-Tribromophenol	10.74	330	210485	91.87	ng	-0.01
44) 2-Fluorobiphenyl	9.27	172	1167195	69.59	ng	-0.01
78) Terphenyl-d14	13.02	244	778976	66.89	ng	-0.01
Target Compounds						
10) Phenol	6.55	94	31688	2.469	ng	Qvalue 91
49) Dimethylphthalate	9.66	163	173985	8.233	ng	100
70) Phenanthrene	11.46	178	144044	5.870	ng	98
74) Fluoranthene	12.65	202	261888	10.539	ng	99
77) Pyrene	12.88	202	256058	12.679	ng	99
80) Benzo(a)anthracene	14.06	228	121870	7.599	ng	96
82) Chrysene	14.10	228	115701	7.578	ng	96
85) Indeno(1,2,3-cd)pyrene	17.07	276	64204	5.257	ng	99
87) Benzo(b)fluoranthene	15.13	252	162755m	10.159	ng	
88) Benzo(k)fluoranthene	15.15	252	49543m	3.131	ng	
89) Benzo(a)pyrene	15.50	252	121511	8.262	ng	# 95
91) Benzo(a,h,i)perylene	17.53	276	60757	5.222	ng	95

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

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