

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092517\
 Data File : BF098821.D
 Acq On : 26 Sep 2017 9:56
 Operator : SJ/JU
 Sample : I5362-05DL 5X
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-2-C-COMPDL

Manual Integrations
 APPROVED

Sohil
 9/26/2017 4:59:07 PM

Quant Time: Sep 26 11:32:04 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	6.91	152	152023	20.00	ng	-0.01
21) Naphthalene-d8	8.19	136	630561	20.00	ng	-0.01
38) Acenaphthene-d10	9.95	164	283184	20.00	ng	-0.01
63) Phenanthrene-d10	11.44	188	465273	20.00	ng	0.00
75) Chrysene-d12	14.07	240	264157	20.00	ng	-0.01
86) Perylene-d12	15.56	264	262820	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	189435	19.84	ng	-0.02
7) Phenol-d6	6.52	99	236589	20.08	ng	-0.02
23) Nitrobenzene-d5	7.47	82	135848	13.82	ng	-0.02
41) 2,4,6-Tribromophenol	10.73	330	40506	17.48	ng	-0.02
44) 2-Fluorobiphenyl	9.26	172	268359	15.82	ng	-0.02
78) Terphenyl-d14	13.02	244	182406	15.70	ng	-0.02
Target Compounds						
31) Naphthalene	8.22	128	76043	2.568	ng	99
37) 2-Methylnaphthalene	8.90	142	41735	2.168	ng	97
51) Acenaphthene	9.98	154	38156	2.153	ng	99
54) Dibenzofuran	10.15	168	54409	2.306	ng	99
70) Phenanthrene	11.46	178	516379	20.734	ng	99
71) Anthracene	11.51	178	104508	4.101	ng	99
74) Fluoranthene	12.65	202	562809	22.317	ng	99
77) Pyrene	12.88	202	466735	23.171	ng	99
80) Benzo(a)anthracene	14.06	228	215955	13.501	ng	98
82) Chrysene	14.10	228	185735	12.197	ng	97
85) Indeno(1,2,3-cd)pyrene	17.06	276	81334	6.677	ng	96
87) Benzo(b)fluoranthene	15.12	252	236483m	14.635	ng	
88) Benzo(k)fluoranthene	15.15	252	58694m	3.677	ng	
89) Benzo(a)pyrene	15.50	252	152192	10.260	ng	95
91) Benzo(a,h,i)perylene	17.52	276	74057	6.311	ng	96

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

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