

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101617\
 Data File : BF099573.D
 Acq On : 17 Oct 2017 4:41
 Operator : SJ/JU
 Sample : I5821-03RE
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WPSB-6A-2-BGSRE

Manual Integrations
 APPROVED

Sohil
 10/17/2017 5:02:13 PM

Quant Time: Oct 17 05:38:06 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 16:25:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	113134	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	448714	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	182605	20.00	ng	-0.01
63) Phenanthrene-d10	11.37	188	263013	20.00	ng	0.00
75) Chrysene-d12	14.02	240	203591	20.00	ng	0.00
86) Perylene-d12	15.49	264	213857	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	759926	106.93	ng	0.00
7) Phenol-d6	6.49	99	908007	103.57	ng	0.00
23) Nitrobenzene-d5	7.41	82	536894	76.73	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	152540	102.09	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	973949	89.04	ng	-0.01
78) Terphenyl-d14	12.95	244	621572	69.43	ng	0.00
Target Compounds						
10) Phenol	6.49	94	55481	5.658	ng	84
49) Dimethylphthalate	9.59	163	167876	12.181	ng	98
70) Phenanthrene	11.39	178	140205	9.959	ng	98
71) Anthracene	11.45	178	32946	2.287	ng	100
74) Fluoranthene	12.59	202	237129	16.634	ng	97
77) Pyrene	12.82	202	221325	14.256	ng	98
80) Benzo(a)anthracene	14.00	228	126726	10.280	ng	99
82) Chrysene	14.04	228	106513	9.075	ng	97
85) Indeno(1,2,3-cd)pyrene	16.97	276	58530	6.234	ng	98
87) Benzo(b)fluoranthene	15.06	252	161705m	12.298	ng	
88) Benzo(k)fluoranthene	15.08	252	49789m	3.834	ng	
89) Benzo(a)pyrene	15.43	252	109787	9.096	ng	97
91) Benzo(a,h,i)perylene	17.42	276	56598	5.928	ng	97

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

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