

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF120117\
 Data File : BF101063.D
 Acq On : 1 Dec 2017 15:17
 Operator : SJ/JU
 Sample : I6610-14
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-10(0-2)-20171127

Manual Integrations
 APPROVED

Sohil
 12/4/2017 11:56:35 AM

Quant Time: Dec 04 00:21:15 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 01 15:52:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	53646	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	210984	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	97895	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	176180	20.00	ng	0.00
75) Chrysene-d12	14.02	240	114588	20.00	ng	0.00
86) Perylene-d12	15.49	264	128080	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	328622	101.22	ng	0.00
7) Phenol-d6	6.49	99	428605	108.74	ng	0.00
23) Nitrobenzene-d5	7.41	82	273662	77.37	ng	0.00
41) 2,4,6-Tribromophenol	10.67	330	75239	63.98	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	536870	75.03	ng	0.00
78) Terphenyl-d14	12.96	244	416878	79.57	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.49	94	15228	3.475	ng	79
31) Naphthalene	8.15	128	24849	2.390	ng	98
48) Acenaphthylene	9.75	152	30893	2.951	ng	98
49) Dimethylphthalate	9.59	163	37051	4.693	ng	99
51) Acenaphthene	9.92	154	29361	4.684	ng	98
54) Dibenzofuran	10.09	168	24067	2.664	ng	# 97
57) Fluorene	10.43	166	41329	5.628	ng	95
70) Phenanthrene	11.40	178	487502	50.247	ng	98
71) Anthracene	11.45	178	126460	12.879	ng	97
72) Carbazole	11.60	167	41905	4.671	ng	97
74) Fluoranthene	12.60	202	649647	60.406	ng	94
77) Pvrene	12.82	202	562675	71.162	ng	98
79) Butylbenzylphthalate	13.43	149	16178	4.641	ng	91
80) Benzo(a)anthracene	14.01	228	272227m	37.627	ng	
82) Chrysene	14.04	228	258344m	38.449	ng	
83) Bis(2-ethylhexyl)phthalate	13.99	149	151590	30.497	ng	98
85) Indeno(1,2,3-cd)pyrene	16.97	276	127497	21.343	ng	# 89
87) Benzo(b)fluoranthene	15.07	252	318759m	41.550	ng	
88) Benzo(k)fluoranthene	15.09	252	110510m	14.621	ng	
89) Benzo(a)pyrene	15.43	252	237058	33.989	ng	97
90) Dibenzo(a,h)anthracene	16.97	278	35041m	5.699	ng	
91) Benzo(g,h,i)perylene	17.43	276	117200	20.226	ng	# 91

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

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