

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110517\
 Data File : BF100231.D
 Acq On : 5 Nov 2017 18:33
 Operator : SJ/JU
 Sample : I6174-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B7(0-10)COMP

Manual Integrations
 APPROVED

Sohil
 11/6/2017 2:33:33 PM

Quant Time: Nov 06 07:12:02 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 03 15:46:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	91159	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	355260	20.00	ng	0.00
38) Acenaphthene-d10	9.79	164	139765	20.00	ng	-0.02
63) Phenanthrene-d10	11.29	188	202414	20.00	ng	-0.01
75) Chrysene-d12	13.94	240	184292	20.00	ng	-0.01
86) Perylene-d12	15.40	264	158822	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	430593	74.16	ng	0.00
7) Phenol-d6	6.42	99	553896	80.45	ng	-0.01
23) Nitrobenzene-d5	7.33	82	299130	54.21	ng	-0.02
41) 2,4,6-Tribromophenol	10.59	330	91176	71.58	ng	-0.02
44) 2-Fluorobiphenyl	9.12	172	607081	67.01	ng	-0.01
78) Terphenyl-d14	12.86	244	415625	49.29	ng	-0.02
Target Compounds						
10) Phenol	6.43	94	24828	3.285	ng	Qvalue # 65
49) Dimethylphthalate	9.52	163	90841	8.207	ng	100
70) Phenanthrene	11.32	178	75789	6.635	ng	97
71) Anthracene	11.37	178	27044	2.332	ng	97
74) Fluoranthene	12.50	202	172937	14.568	ng	98
77) Pyrene	12.74	202	195933	13.982	ng	98
80) Benzo(a)anthracene	13.93	228	113866	9.746	ng	97
82) Chrysene	13.96	228	99201	9.032	ng	97
85) Indeno(1,2,3-cd)pyrene	16.88	276	41564	4.122	ng	94
87) Benzo(b)fluoranthene	14.98	252	117811m	11.747	ng	
88) Benzo(k)fluoranthene	15.00	252	36384m	3.847	ng	
89) Benzo(a)pyrene	15.34	252	83794	9.301	ng	97
91) Benzo(a,h,i)perylene	17.33	276	38910	4.968	ng	97

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

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