

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101217\
 Data File : BF099407.D
 Acq On : 12 Oct 2017 21:00
 Operator : SJ/JU
 Sample : I5729-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 4W-(2-3)

Manual Integrations
 APPROVED

Sohil
 10/13/2017 3:05:50 PM

Quant Time: Oct 13 02:46:06 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	116605	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	460682	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	183377	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	259452	20.00	ng	0.00
75) Chrysene-d12	14.02	240	217233	20.00	ng	0.00
86) Perylene-d12	15.49	264	188469	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	603990	82.46	ng	0.00
7) Phenol-d6	6.49	99	724145	80.14	ng	0.00
23) Nitrobenzene-d5	7.42	82	422322	58.79	ng	0.00
41) 2,4,6-Tribromophenol	10.68	330	107345	71.54	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	793747	72.26	ng	0.00
78) Terphenyl-d14	12.96	244	489973	51.30	ng	0.00
Target Compounds						
10) Phenol	6.50	94	34909	3.454	ng	Qvalue 94
49) Dimethylphthalate	9.60	163	108059	7.808	ng	100
51) Acenaphthene	9.92	154	33502	2.920	ng	99
57) Fluorene	10.43	166	34888	2.841	ng	99
70) Phenanthrene	11.40	178	499442	35.963	ng	99
71) Anthracene	11.45	178	103005	7.249	ng	98
72) Carbazole	11.60	167	42294	3.039	ng	98
73) Di-n-butylphthalate	11.93	149	47467	2.834	ng	99
74) Fluoranthene	12.59	202	541570	38.511	ng	98
77) Pyrene	12.82	202	492071	29.706	ng	99
80) Benzo(a)anthracene	14.00	228	257133	19.548	ng	99
82) Chrysene	14.04	228	232257	18.546	ng	97
85) Indeno(1,2,3-cd)pyrene	16.97	276	77265	7.713	ng	99
87) Benzo(b)fluoranthene	15.06	252	241089m	20.805	ng	
88) Benzo(k)fluoranthene	15.08	252	81483m	7.119	ng	
89) Benzo(a)pyrene	15.43	252	159708	15.015	ng	98
90) Dibenzo(a,h)anthracene	16.97	278	21813	2.562	ng	# 76
91) Benzo(g,h,i)perylene	17.42	276	73860	8.778	ng	91

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

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