

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101217\
 Data File : BF099408.D
 Acq On : 12 Oct 2017 21:28
 Operator : SJ/JU
 Sample : I5729-08 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 10/13/2017 3:05:52 PM

Quant Time: Oct 13 02:54:56 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	114254	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	463467	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	185667	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	270032	20.00	ng	0.00
75) Chrysene-d12	14.02	240	216485	20.00	ng	0.00
86) Perylene-d12	15.49	264	186030	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	147851	20.60	ng	0.00
7) Phenol-d6	6.48	99	184776	20.87	ng	0.00
23) Nitrobenzene-d5	7.41	82	110145	15.24	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	26027	17.13	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	226697	20.38	ng	0.00
78) Terphenyl-d14	12.95	244	142540	14.97	ng	0.00
Target Compounds						
37) 2-Methylnaphthalene	8.85	142	28513	2.015	ng	96
70) Phenanthrene	11.40	178	222890	15.421	ng	98
71) Anthracene	11.45	178	47218	3.193	ng	97
73) Di-n-butylphthalate	11.93	149	51344	2.945	ng	# 99
74) Fluoranthene	12.59	202	406863	27.798	ng	97
77) Pyrene	12.82	202	376048	22.780	ng	99
80) Benzo(a)anthracene	14.00	228	216819	16.540	ng	99
82) Chrysene	14.04	228	201551	16.150	ng	97
85) Indeno(1,2,3-cd)pyrene	16.97	276	73715	7.384	ng	98
87) Benzo(b)fluoranthene	15.06	252	224165m	19.598	ng	
88) Benzo(k)fluoranthene	15.08	252	74413m	6.587	ng	
89) Benzo(a)pyrene	15.43	252	150866	14.369	ng	# 97
90) Dibenzo(a,h)anthracene	16.97	278	20804m	2.476	ng	
91) Benzo(a,h,i)perylene	17.42	276	70207	8.453	ng	93

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

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