

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101217\
 Data File : BF099411.D
 Acq On : 12 Oct 2017 22:52
 Operator : SJ/JU
 Sample : I5729-11
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 6W-(2-3)

Manual Integrations
 APPROVED

Sohil
 10/13/2017 3:06:28 PM

Quant Time: Oct 13 04:13:45 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	111286	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	442317	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	180805	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	266344	20.00	ng	0.00
75) Chrysene-d12	14.02	240	214508	20.00	ng	0.00
86) Perylene-d12	15.49	264	175063	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	609542	87.19	ng	0.00
7) Phenol-d6	6.49	99	729320	84.57	ng	0.00
23) Nitrobenzene-d5	7.42	82	428891	62.18	ng	0.00
41) 2,4,6-Tribromophenol	10.68	330	112754	76.21	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	755634	69.77	ng	0.00
78) Terphenyl-d14	12.96	244	484094	51.32	ng	0.00
Target Compounds						
48) Acenaphthylene	9.75	152	56000	3.135	ng	99
49) Dimethylphthalate	9.60	163	101053	7.405	ng	100
70) Phenanthrene	11.40	178	79336	5.565	ng	97
71) Anthracene	11.45	178	41578	2.850	ng	97
74) Fluoranthene	12.59	202	122988	8.519	ng	98
77) Pyrene	12.82	202	108671	6.644	ng	97
80) Benzo(a)anthracene	14.00	228	58404m	4.496	ng	
82) Chrysene	14.04	228	53472	4.324	ng	97
85) Indeno(1,2,3-cd)pyrene	16.96	276	41336	4.179	ng	99
87) Benzo(b)fluoranthene	15.06	252	79225m	7.360	ng	
89) Benzo(a)pyrene	15.42	252	55816	5.649	ng	# 93
91) Benzo(a,h,i)perylene	17.42	276	47726	6.106	ng	# 90

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

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