

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101217\
 Data File : BF099404.D
 Acq On : 12 Oct 2017 19:36
 Operator : SJ/JU
 Sample : I5729-05 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 3E-(1-2)

Manual Integrations
 APPROVED

Sohil
 10/13/2017 3:05:44 PM

Quant Time: Oct 13 02:21:54 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	112169	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	432979	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	175084	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	253626	20.00	ng	0.00
75) Chrysene-d12	14.02	240	206588	20.00	ng	0.00
86) Perylene-d12	15.49	264	182395	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	135501	19.23	ng	0.00
7) Phenol-d6	6.48	99	164748	18.95	ng	0.00
23) Nitrobenzene-d5	7.41	82	91577	13.56	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	23098	16.12	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	167776	16.00	ng	0.00
78) Terphenyl-d14	12.95	244	99273	10.93	ng	0.00
Target Compounds						Qvalue
51) Acenaphthene	9.92	154	29641	2.706	ng	100
54) Dibenzofuran	10.09	168	34521	2.367	ng	98
57) Fluorene	10.43	166	37108	3.165	ng	97
70) Phenanthrene	11.40	178	389113	28.662	ng	98
71) Anthracene	11.45	178	87942	6.331	ng	99
72) Carbazole	11.61	167	45767	3.365	ng	99
74) Fluoranthene	12.59	202	610735	44.426	ng	99
77) Pyrene	12.82	202	597933	37.956	ng	99
80) Benzo(a)anthracene	14.01	228	275455	22.020	ng	99
82) Chrysene	14.04	228	264582	22.216	ng	98
85) Indeno(1,2,3-cd)pyrene	16.97	276	96501	10.129	ng	98
87) Benzo(b)fluoranthene	15.06	252	281573m	25.108	ng	
88) Benzo(k)fluoranthene	15.09	252	104156m	9.403	ng	
89) Benzo(a)pyrene	15.43	252	176682	17.164	ng	# 94
91) Benzo(a,h,i)perylene	17.42	276	92238	11.327	ng	92

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

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