

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111117\
 Data File : BF100442.D
 Acq On : 11 Nov 2017 15:56
 Operator : SJ/JU
 Sample : I6399-22
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-15

Manual Integrations
 APPROVED

SOHIL
 11/13/2017 4:54:25 PM

Quant Time: Nov 13 03:29:06 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 10 15:44:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	149239	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	527475	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	194304	20.00	ng	0.00
63) Phenanthrene-d10	11.43	188	315928	20.00	ng	0.00
75) Chrysene-d12	14.07	240	317582	20.00	ng	0.00
86) Perylene-d12	15.55	264	230414	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	808162	86.81	ng	0.01
7) Phenol-d6	6.53	99	967445	84.27	ng	0.00
23) Nitrobenzene-d5	7.46	82	574069	71.95	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	161638	81.64	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	886282	69.46	ng	0.00
78) Terphenyl-d14	13.01	244	666604	48.23	ng	0.00
Target Compounds						
10) Phenol	6.54	94	50940	4.227	ng	97
49) Dimethylphthalate	9.65	163	87049	5.868	ng	99
74) Fluoranthene	12.64	202	122106	6.926	ng	95
77) Pyrene	12.87	202	135162	5.970	ng	98
80) Benzo(a)anthracene	14.06	228	41810	2.186	ng	97
82) Chrysene	14.09	228	47285	2.553	ng	98
87) Benzo(b)fluoranthene	15.12	252	54116m	3.964	ng	
89) Benzo(a)pyrene	15.49	252	32462	2.682	ng	98
91) Benzo(g,h,i)perylene	17.50	276	24141	2.492	ng	# 89

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

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