

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101517\
 Data File : BF099513.D
 Acq On : 15 Oct 2017 14:57
 Operator : SJ/JU
 Sample : I5752-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ENV-118-7(0-5)

Manual Integrations
 APPROVED

Sohil
 10/16/2017 7:45:25 PM

Quant Time: Oct 16 01:55:34 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	106433	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	428668	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	174506	20.00	ng	-0.01
63) Phenanthrene-d10	11.37	188	255365	20.00	ng	0.00
75) Chrysene-d12	14.01	240	186456	20.00	ng	0.00
86) Perylene-d12	15.49	264	209683	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	767623	114.81	ng	0.00
7) Phenol-d6	6.49	99	916728	111.15	ng	0.00
23) Nitrobenzene-d5	7.41	82	542262	81.12	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	147388	103.22	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	891896	85.32	ng	-0.01
78) Terphenyl-d14	12.95	244	518494	63.24	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.49	94	57050	6.185	ng	95
49) Dimethylphthalate	9.59	163	307630	23.357	ng	100
70) Phenanthrene	11.39	178	153086	11.200	ng	98
71) Anthracene	11.45	178	36413	2.604	ng	99
74) Fluoranthene	12.58	202	124991	9.030	ng	99
77) Pyrene	12.82	202	110589	7.778	ng	97
80) Benzo(a)anthracene	14.00	228	55752	4.938	ng	96
82) Chrysene	14.04	228	49322	4.589	ng	97
85) Indeno(1,2,3-cd)pyrene	16.96	276	29410	3.420	ng	97
87) Benzo(b)fluoranthene	15.05	252	60460m	4.690	ng	
88) Benzo(k)fluoranthene	15.08	252	25789m	2.025	ng	
89) Benzo(a)pyrene	15.42	252	49545	4.187	ng	# 96
91) Benzo(a,h,i)perylene	17.42	276	27096	2.894	ng	92

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

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