

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090617\
 Data File : BF098282.D
 Acq On : 6 Sep 2017 19:42
 Operator : SJ/JU
 Sample : I5095-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 STORM-SEWER-SED

Manual Integrations
 APPROVED

Sohil
 9/7/2017 5:44:32 PM

Quant Time: Sep 07 05:36:06 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 17:28:20 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	103340	20.00	ng	-0.01
21) Naphthalene-d8	7.97	136	402202	20.00	ng	-0.01
38) Acenaphthene-d10	9.73	164	171912	20.00	ng	-0.01
63) Phenanthrene-d10	11.21	188	279779	20.00	ng	-0.01
75) Chrysene-d12	13.85	240	200306	20.00	ng	-0.01
86) Perylene-d12	15.26	264	179945	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	617633	90.91	ng	0.00
7) Phenol-d6	6.34	99	844588	91.50	ng	-0.01
23) Nitrobenzene-d5	7.26	82	500465	67.60	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	117590	78.83	ng	-0.01
44) 2-Fluorobiphenyl	9.05	172	579519	49.53	ng	-0.02
78) Terphenyl-d14	12.80	244	339327	35.33	ng	-0.01
Target Compounds						
31) Naphthalene	8.00	128	51643	2.473	ng	Qvalue 98
48) Acenaphthylene	9.59	152	117503	6.460	ng	99
49) Dimethylphthalate	9.45	163	125414	9.980	ng	99
51) Acenaphthene	9.76	154	47338	4.126	ng	98
54) Dibenzofuran	9.93	168	36824	2.395	ng	# 95
57) Fluorene	10.27	166	77753	6.541	ng	97
70) Phenanthrene	11.24	178	824817	55.325	ng	99
71) Anthracene	11.29	178	261237	17.276	ng	99
72) Carbazole	11.45	167	88228	6.147	ng	99
73) Di-n-butylphthalate	11.78	149	37993	2.298	ng	97
74) Fluoranthene	12.43	202	1467239	94.289	ng	99
77) Pvrene	12.66	202	1281752	78.238	ng	99
79) Butylbenzylphthalate	13.27	149	29121	4.636	ng	# 69
80) Benzo(a)anthracene	13.85	228	763241	63.576	ng	97
82) Chrysene	13.88	228	725172	60.723	ng	97
83) Bis(2-ethylhexyl)phthalate	13.83	149	100302	14.411	ng	# 97
85) Indeno(1,2,3-cd)pvrene	16.63	276	312768m	35.602	ng	
87) Benzo(b)fluoranthene	14.87	252	896030m	78.998	ng	
88) Benzo(k)fluoranthene	14.89	252	258484m	24.256	ng	
89) Benzo(a)pvrene	15.22	252	595550	59.310	ng	98
90) Dibenzo(a,h)anthracene	16.63	278	81271	9.942	ng	# 74
91) Benzo(g,h,i)perylene	17.04	276	279069	32.717	ng	95

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

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