

Data Path : Z:\HPCHEM1\BNA F\DATA\BF103017\
 Data File : BF099967.D
 Acq On : 30 Oct 2017 18:37
 Operator : SJ/JU
 Sample : I6171-09 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-102617-E

Manual Integrations
 APPROVED

Sohil
 10/31/2017 7:41:10 PM

Quant Time: Oct 31 02:59:41 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 27 15:58:57 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	96686	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	380510	20.00	ng	-0.01
38) Acenaphthene-d10	9.83	164	164833	20.00	ng	-0.01
63) Phenanthrene-d10	11.32	188	265759	20.00	ng	-0.01
75) Chrysene-d12	13.99	240	169744	20.00	ng	0.00
86) Perylene-d12	15.47	264	162446	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	334435	54.30	ng	-0.01
7) Phenol-d6	6.43	99	419348	57.43	ng	-0.01
23) Nitrobenzene-d5	7.36	82	233317	39.48	ng	-0.01
41) 2,4,6-Tribromophenol	10.62	330	84265	56.10	ng	-0.02
44) 2-Fluorobiphenyl	9.15	172	504941	47.26	ng	-0.01
78) Terphenyl-d14	12.91	244	356193	45.86	ng	-0.01
Target Compounds						
10) Phenol	6.45	94	22785	2.842	ng	Qvalue 85
31) Naphthalene	8.09	128	67270	3.662	ng	97
37) 2-Methylnaphthalene	8.79	142	51792	4.208	ng	96
48) Acenaphthylene	9.69	152	192259	11.527	ng	99
49) Dimethylphthalate	9.54	163	86639	6.637	ng	99
51) Acenaphthene	9.86	154	64077	6.287	ng	99
54) Dibenzofuran	10.03	168	79149	5.502	ng	97
55) 4-Nitrophenol	10.03	139	28332	16.770	ng	# 12
57) Fluorene	10.37	166	144390	12.122	ng	99
70) Phenanthrene	11.35	178	918550	61.245	ng	98
71) Anthracene	11.40	178	350482	23.022	ng	99
72) Carbazole	11.56	167	80791	5.643	ng	98
74) Fluoranthene	12.55	202	1352107	86.749	ng	99
77) Pyrene	12.78	202	1214827	94.120	ng	100
80) Benzo(a)anthracene	13.97	228	653407	60.722	ng	99
82) Chrysene	14.02	228	545822	53.954	ng	98
85) Indeno(1,2,3-cd)pyrene	16.97	276	302959	32.622	ng	95
87) Benzo(b)fluoranthene	15.05	252	814323m	79.387	ng	
88) Benzo(k)fluoranthene	15.07	252	213799m	22.103	ng	
89) Benzo(a)pyrene	15.42	252	542954	58.925	ng	97
90) Dibenzo(a,h)anthracene	16.97	278	66827m	8.162	ng	
91) Benzo(g,h,i)perylene	17.43	276	265031	33.086	ng	96

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

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