

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
 Data File : BF101780.D
 Acq On : 27 Dec 2017 23:16
 Operator : SJ/JU
 Sample : I7064-07 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAMPLE-4

Manual Integrations
 APPROVED

Sohil
 12/28/2017 2:03:08 PM

Quant Time: Dec 28 05:52:09 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 10:48:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	155331	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	588623	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	233788	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	369054	20.00	ng	0.00
75) Chrysene-d12	13.97	240	296948	20.00	ng	0.00
86) Perylene-d12	15.44	264	264030	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	586056	56.20	ng	0.01
7) Phenol-d6	6.43	99	701836	54.08	ng	0.00
23) Nitrobenzene-d5	7.36	82	414073	39.16	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	116883	51.89	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	696572	46.22	ng	0.00
78) Terphenyl-d14	12.90	244	474332	38.51	ng	0.00
Target Compounds						
10) Phenol	6.46	94	33224	2.246	ng	Qvalue # 66
49) Dimethylphthalate	9.55	163	62454	3.321	ng	99
51) Acenaphthene	9.86	154	33716	2.240	ng	99
57) Fluorene	10.38	166	43462	2.685	ng	98
70) Phenanthrene	11.35	178	475969	23.191	ng	98
71) Anthracene	11.40	178	148237	7.071	ng	98
74) Fluoranthene	12.54	202	599473	27.827	ng	99
77) Pyrene	12.77	202	516316	23.168	ng	99
80) Benzo(a)anthracene	13.96	228	298687	15.233	ng	98
82) Chrysene	14.00	228	258357	13.955	ng	99
85) Indeno(1,2,3-cd)pyrene	16.94	276	98783	5.992	ng	# 91
87) Benzo(b)fluoranthene	15.02	252	268562m	16.688	ng	
88) Benzo(k)fluoranthene	15.04	252	113165m	7.232	ng	
89) Benzo(a)pyrene	15.39	252	189335	12.762	ng	93
91) Benzo(a,h,i)perylene	17.39	276	90966	7.212	ng	94

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

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