

Data Path : Z:\HPCHEM1\BNA F\DATA\BF021318\
 Data File : BF102973.D
 Acq On : 13 Feb 2018 20:28
 Operator : SJ/JU
 Sample : J1537-06
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-03-COMP

Manual Integrations
 APPROVED

Sohil
 2/14/2018 10:29:49 AM

Quant Time: Feb 14 02:36:47 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 00:38:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	155369	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	610646	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	229589	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	351563	20.00	ng	0.00
75) Chrysene-d12	14.04	240	240041	20.00	ng	0.00
86) Perylene-d12	15.53	264	176738	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1253742	122.55	ng	0.01
7) Phenol-d6	6.51	99	1466568	119.06	ng	0.01
23) Nitrobenzene-d5	7.44	82	904926	102.80	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	235781	127.59	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1276485	92.26	ng	0.00
78) Terphenyl-d14	12.99	244	930967	91.83	ng	0.00
Target Compounds						
10) Phenol	6.52	94	107249	8.124	ng	98
49) Dimethylphthalate	9.62	163	100001	5.586	ng	100
51) Acenaphthene	9.95	154	30365	2.147	ng	95
57) Fluorene	10.46	166	37082	2.558	ng	98
70) Phenanthrene	11.42	178	143723	7.340	ng	100
71) Anthracene	11.47	178	54275	2.725	ng	97
74) Fluoranthene	12.62	202	224557	11.399	ng	98
77) Pyrene	12.84	202	266915	14.180	ng	98
80) Benzo(a)anthracene	14.03	228	98079	6.497	ng	# 92
82) Chrysene	14.07	228	102734	6.751	ng	# 94
85) Indeno(1,2,3-cd)pyrene	17.02	276	45644	3.616	ng	97
87) Benzo(b)fluoranthene	15.09	252	101366m	8.846	ng	
88) Benzo(k)fluoranthene	15.12	252	33534m	3.063	ng	
89) Benzo(a)pyrene	15.46	252	78087	7.571	ng	# 71
91) Benzo(a,h,i)perylene	17.48	276	46984	5.734	ng	# 87

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

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