

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030218\
 Data File : BF103389.D
 Acq On : 3 Mar 2018 1:53
 Operator : SJ/JU
 Sample : J1724-11
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-19

Manual Integrations
 APPROVED

Sohil
 3/5/2018 3:14:14 PM

Quant Time: Mar 05 04:29:27 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 02 13:28:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	141275	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	600546	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	238796	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	341494	20.00	ng	0.00
75) Chrysene-d12	14.07	240	205786	20.00	ng	0.00
86) Perylene-d12	15.58	264	148018	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	943969	101.06	ng	0.00
7) Phenol-d6	6.52	99	1131162	98.96	ng	0.00
23) Nitrobenzene-d5	7.45	82	719009	68.37	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	154081	76.23	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	965894	66.01	ng	0.00
78) Terphenyl-d14	13.00	244	599622	70.04	ng	0.00
Target Compounds						
10) Phenol	6.53	94	32185	2.426	ng	Qvalue # 62
49) Dimethylphthalate	9.63	163	79203	4.135	ng	99
70) Phenanthrene	11.44	178	273601	14.558	ng	99
71) Anthracene	11.49	178	88595	4.597	ng	98
74) Fluoranthene	12.63	202	526524	27.880	ng	100
77) Pyrene	12.86	202	473522	29.513	ng	99
80) Benzo(a)anthracene	14.06	228	221306	16.575	ng	98
82) Chrysene	14.09	228	201668	15.555	ng	99
85) Indeno(1,2,3-cd)pyrene	17.12	276	70809	6.899	ng	96
87) Benzo(b)fluoranthene	15.13	252	187544m	19.271	ng	
88) Benzo(k)fluoranthene	15.16	252	74018m	8.077	ng	
89) Benzo(a)pyrene	15.52	252	126926	14.605	ng	# 93
90) Dibenzo(a,h)anthracene	17.12	278	19248	2.866	ng	# 70
91) Benzo(a,h,i)perylene	17.59	276	65979	10.053	ng	# 88

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

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