

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103468.D
 Acq On : 6 Mar 2018 21:02
 Operator : SJ/JU
 Sample : J1722-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-7

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:17:23 PM

Quant Time: Mar 07 01:09:52 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 06 11:52:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	129346	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	533486	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	213567	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	310581	20.00	ng	0.00
75) Chrysene-d12	14.06	240	205740	20.00	ng	0.00
86) Perylene-d12	15.57	264	154990	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	955782	111.76	ng	0.00
7) Phenol-d6	6.51	99	1151357	110.02	ng	0.00
23) Nitrobenzene-d5	7.44	82	730881	78.24	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	135555	74.99	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	954967	74.85	ng	0.00
78) Terphenyl-d14	12.99	244	610153	71.29	ng	0.00
Target Compounds						
10) Phenol	6.52	94	88238	7.263	ng	Qvalue # 72
49) Dimethylphthalate	9.62	163	91845	5.361	ng	100
70) Phenanthrene	11.43	178	127052	7.433	ng	98
71) Anthracene	11.49	178	39697	2.265	ng	97
74) Fluoranthene	12.63	202	247598	14.416	ng	99
77) Pyrene	12.86	202	243125	15.157	ng	99
80) Benzo(a)anthracene	14.05	228	107250	8.034	ng	97
82) Chrysene	14.09	228	102200	7.885	ng	97
83) Bis(2-ethylhexyl)phthalate	14.01	149	25957	2.270	ng	99
85) Indeno(1,2,3-cd)pyrene	17.13	276	39286	3.828	ng	97
87) Benzo(b)fluoranthene	15.13	252	103775m	10.184	ng	
88) Benzo(k)fluoranthene	15.16	252	37733m	3.932	ng	
89) Benzo(a)pyrene	15.52	252	70533	7.751	ng	# 91
91) Benzo(a,h,i)perylene	17.60	276	38790	5.644	ng	# 87

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

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