

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032718\
 Data File : BF103983.D
 Acq On : 27 Mar 2018 21:14
 Operator : JU/SJ
 Sample : J2052-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-BD

Manual Integrations
 APPROVED

Sohil
 3/28/2018 4:54:28 PM

Quant Time: Mar 28 05:49:34 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 23 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	131683	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	527766	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	226620	20.00	ng	-0.01
63) Phenanthrene-d10	11.50	188	355423	20.00	ng	0.00
75) Chrysene-d12	14.16	240	249002	20.00	ng	-0.01
86) Perylene-d12	15.70	264	214166	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	744396	85.98	ng	0.01
7) Phenol-d6	6.60	99	912584	86.16	ng	0.00
23) Nitrobenzene-d5	7.53	82	526973	69.63	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	199850	102.57	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	903052	63.56	ng	0.00
78) Terphenyl-d14	13.09	244	677626	63.17	ng	-0.01
Target Compounds						
10) Phenol	6.62	94	23202	2.061	ng	Qvalue # 57
48) Acenaphthylene	9.87	152	51404	2.209	ng	97
49) Dimethylphthalate	9.72	163	67994	3.933	ng	99
70) Phenanthrene	11.53	178	102835	5.289	ng	96
71) Anthracene	11.58	178	44712	2.264	ng	98
74) Fluoranthene	12.72	202	385443	19.243	ng	99
77) Pyrene	12.96	202	385597	21.086	ng	98
80) Benzo(a)anthracene	14.14	228	196803	12.486	ng	98
82) Chrysene	14.19	228	172267	11.465	ng	97
85) Indeno(1,2,3-cd)pyrene	17.28	276	85090	6.485	ng	# 93
87) Benzo(b)fluoranthene	15.24	252	288123m	21.294	ng	
88) Benzo(k)fluoranthene	15.27	252	65238m	5.016	ng	
89) Benzo(a)pyrene	15.63	252	150748	12.284	ng	96
90) Dibenzo(a,h)anthracene	17.29	278	28141m	2.648	ng	
91) Benzo(a,h,i)perylene	17.76	276	69913	6.763	ng	93

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

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