

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
 Data File : BF113494.D
 Acq On : 2 Apr 2019 6:37
 Operator : JU/SJ
 Sample : K2056-03DL 2X
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-9DL

Manual Integrations
 APPROVED

Sohil
 4/2/2019 3:49:57 PM

Quant Time: Apr 02 07:58:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	132528	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	499839	20.00	ng	0.00
39) Acenaphthene-d10	9.74	164	219409	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	395006	20.00	ng	0.00
76) Chrysene-d12	13.85	240	363840	20.00	ng	0.00
87) Perylene-d12	15.26	264	249175	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	374693	46.23	ng	0.00
7) Phenol-d6	6.35	99	517008	50.42	ng	0.00
23) Nitrobenzene-d5	7.27	82	266983	31.70	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	116429	42.96	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	581831	30.19	ng	0.00
79) Terphenyl-d14	12.80	244	574120	34.78	ng	0.00
Target Compounds						
49) Acenaphthylene	9.60	152	49721	2.179	ng	99
52) Acenaphthene	9.77	154	34868	2.714	ng	99
58) Fluorene	10.29	166	31473	2.098	ng	97
71) Phenanthrene	11.24	178	340685	17.368	ng	98
72) Anthracene	11.29	178	128865	6.468	ng	98
73) Carbazole	11.45	167	82389	4.333	ng	98
75) Fluoranthene	12.43	202	1594636	71.909	ng	100
78) Pyrene	12.66	202	1362697	54.048	ng	99
81) Benzo(a)anthracene	13.84	228	386469	18.556	ng	97
83) Chrysene	13.87	228	819901	38.758	ng	99
86) Indeno(1,2,3-cd)pyrene	16.62	276	78592	4.329	ng	95
88) Benzo(b)fluoranthene	14.86	252	549785m	35.519	ng	
89) Benzo(k)fluoranthene	14.88	252	178198m	13.109	ng	
90) Benzo(a)pyrene	15.20	252	120275	8.765	ng	97
92) Benzo(g,h,i)perylene	17.03	276	62067	5.188	ng	93

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

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