

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP092719\
 Data File : BP000466.D
 Acq On : 27 Sep 2019 18:44
 Operator : JU
 Sample : K5038-03DL 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DN-2DL

Manual Integrations
 APPROVED

mohammad
 9/30/2019 8:15:28 AM

Quant Time: Sep 28 00:05:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 19 14:44:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	218871	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	376695	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	209409	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	386654	20.00	ng	0.00
76) Chrysene-d12	14.00	240	640672	20.00	ng	0.00
87) Perylene-d12	15.49	264	388285	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	236285	18.64	ng	0.00
7) Phenol-d6	6.42	99	352749	22.70	ng	0.00
23) Nitrobenzene-d5	7.34	82	120589	20.65	ng	-0.01
42) 2,4,6-Tribromophenol	10.63	330	53161	25.59	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	218808	18.81	ng	0.00
79) Terphenyl-d14	12.93	244	537370	17.60	ng	0.00
Target Compounds						
49) Acenaphthylene	9.69	152	38018	2.081	ng	99
50) Dimethylphthalate	9.55	163	47437	3.325	ng	99
71) Phenanthrene	11.35	178	148217	7.649	ng	99
72) Anthracene	11.40	178	99815	5.005	ng	98
75) Fluoranthene	12.56	202	1058742	50.887	ng	98
78) Pyrene	12.79	202	996657	20.793	ng	99
81) Benzo(a)anthracene	13.99	228	579900	14.329	ng	97
83) Chrysene	14.03	228	536944	13.513	ng	97
86) Indeno(1,2,3-cd)pyrene	16.97	276	191343	3.951	ng	# 87
88) Benzo(b)fluoranthene	15.06	252	408765m	17.647	ng	
89) Benzo(k)fluoranthene	15.09	252	150099m	7.429	ng	
90) Benzo(a)pyrene	15.43	252	274413m	13.020	ng	
91) Dibenzo(a,h)anthracene	16.98	278	47789	2.432	ng	# 92
92) Benzo(g,h,i)perylene	17.42	276	174808	8.654	ng	# 91

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

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