

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091420\
 Data File : BP003283.D
 Acq On : 14 Sep 2020 23:28
 Operator : CG/JU
 Sample : L3981-09 10X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 B19-7-9

Manual Integrations
 APPROVED

mohammad
 9/15/2020 1:22:14 PM

Quant Time: Sep 15 05:18:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 01 14:37:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	181049	20.00	ng	-0.01
21) Naphthalene-d8	10.42	136	680579	20.00	ng	-0.01
39) Acenaphthene-d10	14.29	164	385293	20.00	ng	0.00
64) Phenanthrene-d10	17.04	188	694625	20.00	ng	0.00
76) Chrysene-d12	21.14	240	683845	20.00	ng	0.00
86) Perylene-d12	23.37	264	832292	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	69305	6.77	ng	0.00
7) Phenol-d6	6.84	99	83974	6.54	ng	0.00
23) Nitrobenzene-d5	8.79	82	49632	5.26	ng	0.00
42) 2,4,6-Tribromophenol	15.79	330	34065	6.54	ng	0.00
45) 2-Fluorobiphenyl	12.90	172	146211	5.56	ng	-0.01
79) Terphenyl-d14	19.62	244	151837	4.89	ng	0.00

Target Compounds				Qvalue		
31) Naphthalene	10.47	128	191092	5.465	ng	99
37) 2-Methylnaphthalene	12.09	142	83121	3.323	ng	99
38) 1-Methylnaphthalene	12.32	142	81269	3.418	ng	96
49) Acenaphthylene	14.00	152	352950	9.867	ng	99
52) Acenaphthene	14.35	154	81879	3.726	ng	99
55) Dibenzofuran	14.69	168	152181	4.409	ng	97
58) Fluorene	15.34	166	189324	7.121	ng	99
71) Phenanthrene	17.09	178	2626874	69.791	ng	99
72) Anthracene	17.17	178	421290	11.695	ng	99
73) Carbazole	17.45	167	164631	4.762	ng	98
75) Fluoranthene	19.07	202	3037630	69.827	ng	98
78) Pyrene	19.42	202	3229456	72.763	ng	99
81) Benzo(a)anthracene	21.13	228	1474137	34.414	ng	97
83) Chrysene	21.17	228	1591975	38.816	ng	98
87) Indeno(1,2,3-cd)pyrene	25.66	276	1011511	16.297	ng	# 94
88) Benzo(b)fluoranthene	22.71	252	1852919m	35.819	ng	
89) Benzo(k)fluoranthene	22.75	252	672410m	13.605	ng	
90) Benzo(a)pyrene	23.28	252	1519712	31.667	ng	97
91) Dibenzo(a,h)anthracene	25.65	278	270936m	5.314	ng	
92) Benzo(a,h,i)perylene	26.35	276	1022425	19.981	ng	96

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

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